

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007652.D
 Acq On : 25 Oct 2021 21:39
 Operator : CG/JU
 Sample : PB140192BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK192

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

Signal : TIC: BP007652.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.370	10	14	20	rBV	53165	66279	1.34%	0.141%
2	3.781	81	84	101	rBV	665798	1039067	21.08%	2.212%
3	4.122	138	142	146	rVB2	12057	15615	0.32%	0.033%
4	7.105	642	649	660	rBV	751488	1181141	23.96%	2.515%
5	7.269	669	677	688	rVB	589130	942352	19.12%	2.006%
6	7.469	703	711	722	rBV	785068	1226747	24.89%	2.612%
7	7.940	784	791	798	rBV	579223	924304	18.75%	1.968%
8	8.646	903	911	927	rBV	930664	1487349	30.18%	3.166%
9	9.093	980	987	995	rBV	682812	1125243	22.83%	2.396%
10	9.822	1103	1111	1118	rBV	554798	906452	18.39%	1.930%
11	10.363	1195	1203	1214	rBV	889728	1495650	30.35%	3.184%
12	10.740	1260	1267	1279	rVB	765278	1316704	26.72%	2.803%
13	10.875	1279	1290	1303	rBV	864425	1518295	30.81%	3.232%
14	12.328	1533	1537	1544	rVB2	17382	30974	0.63%	0.066%
15	13.545	1741	1744	1752	rVB9	4519	7395	0.15%	0.016%
16	13.987	1812	1819	1832	rBV	1659608	2291723	46.50%	4.879%
17	14.269	1857	1867	1879	rBV	1736041	2640757	53.58%	5.622%
18	14.575	1911	1919	1929	rBV	1198993	1797462	36.47%	3.827%
19	14.763	1945	1951	1964	rBV	732933	1093115	22.18%	2.327%
20	15.569	2081	2088	2097	rBV	2344189	3357823	68.13%	7.149%
21	15.681	2101	2107	2120	rBV	630587	870280	17.66%	1.853%
22	15.810	2125	2129	2135	rVB	16677	30339	0.62%	0.065%
23	16.251	2198	2204	2210	rVB	5346	9795	0.20%	0.021%
24	17.010	2329	2333	2338	rBV7	3871	5571	0.11%	0.012%
25	17.116	2343	2351	2357	rBV2	8335	16920	0.34%	0.036%
26	17.328	2380	2387	2394	rBV	1546017	2213149	44.90%	4.712%
27	17.428	2397	2404	2417	rBV2	2572826	3761443	76.32%	8.008%
28	17.704	2450	2451	2459	rVB7	4276	5484	0.11%	0.012%
29	18.228	2535	2540	2542	rBV6	5563	8996	0.18%	0.019%
30	18.728	2621	2625	2628	rBV6	5843	8370	0.17%	0.018%
31	19.239	2708	2712	2716	rBV2	40123	54349	1.10%	0.116%
32	19.304	2719	2723	2727	rBV	39983	55879	1.13%	0.119%
33	19.663	2778	2784	2795	rBV	3493888	4567168	92.67%	9.723%
34	21.151	3033	3037	3040	rBV2	62952	90945	1.85%	0.194%
35	21.192	3041	3044	3051	rVV	217703	291368	5.91%	0.620%
36	21.416	3077	3082	3088	rBV	2039506	2720710	55.20%	5.792%

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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

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37	23.704	3464	3471	3480	rVB2	2500928	4928663	100.00%	10.493%
38	23.857	3491	3497	3508	rVB2	1344095	2867772	58.19%	6.105%

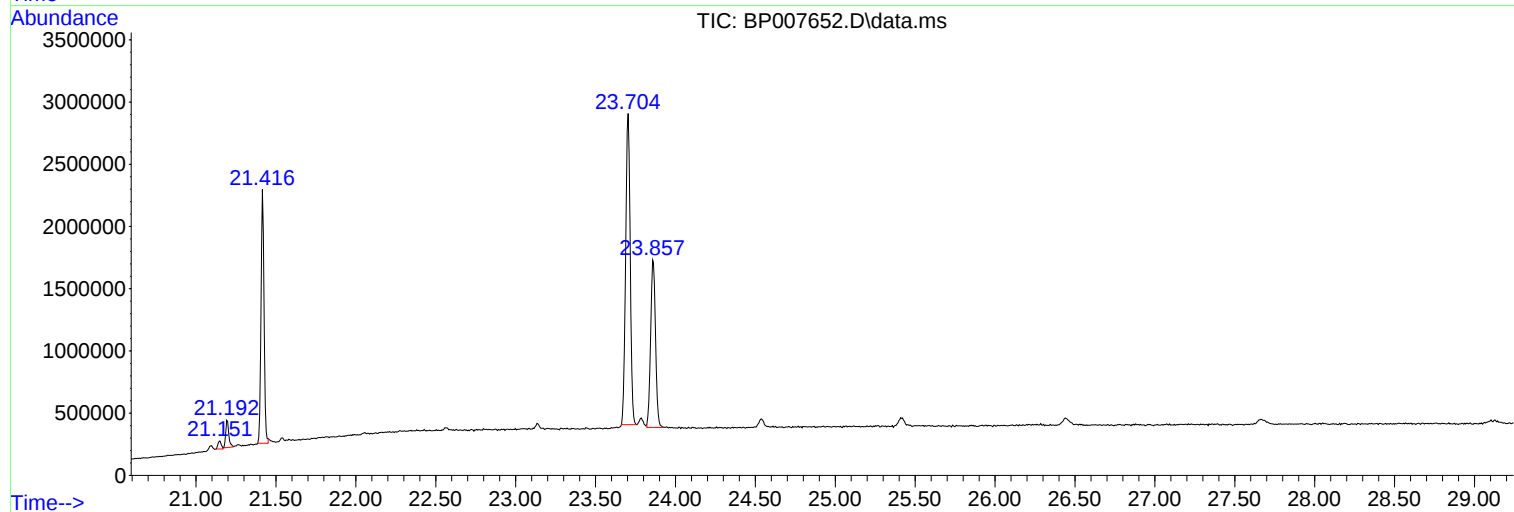
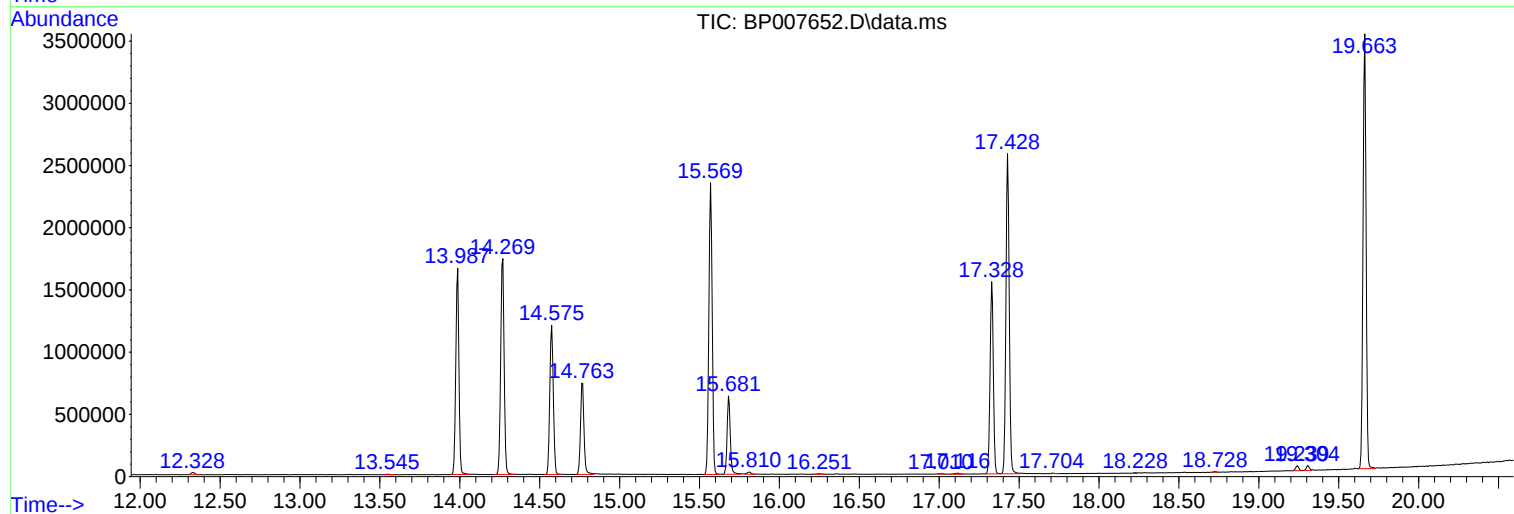
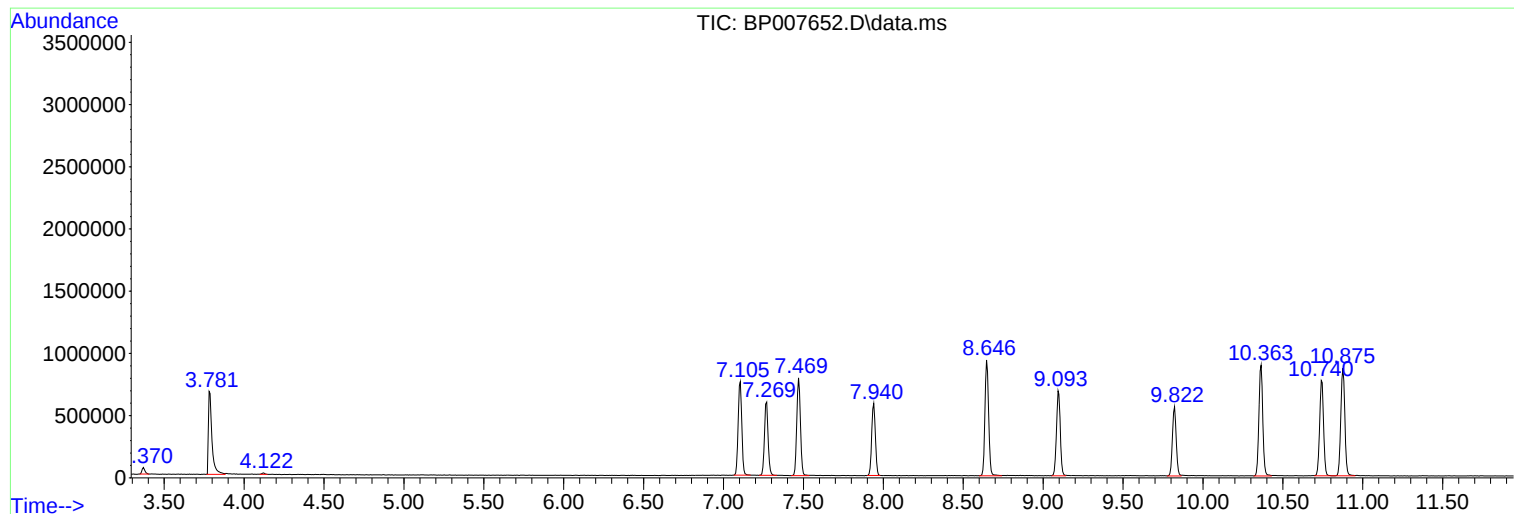
Sum of corrected areas: 46971648

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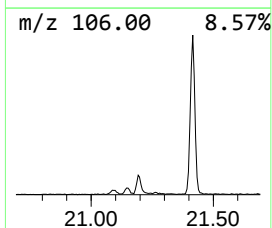
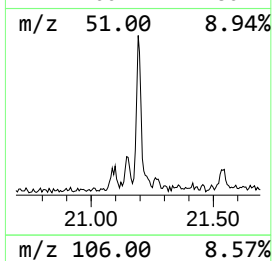
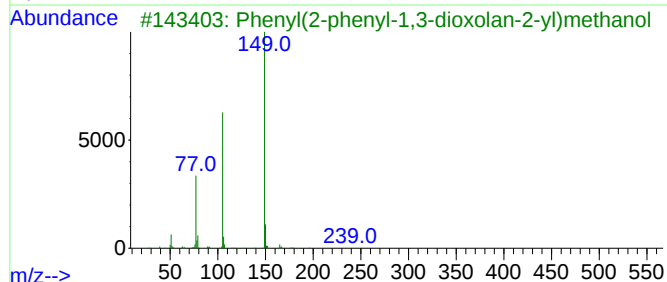
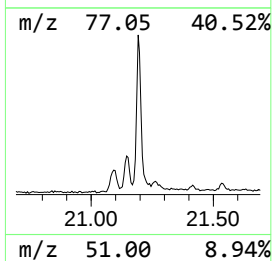
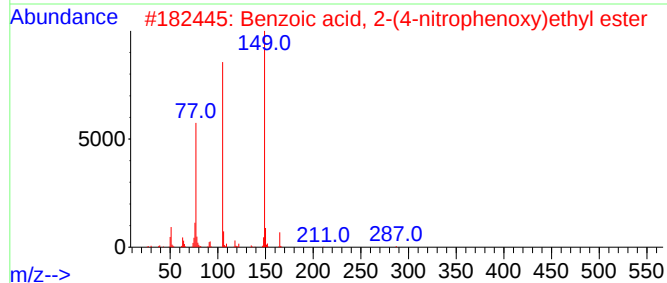
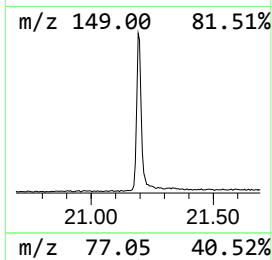
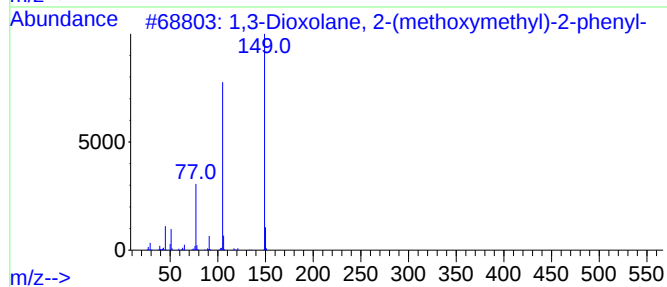
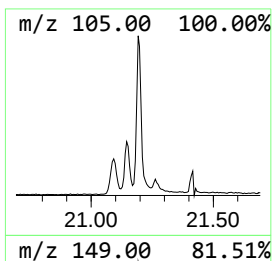
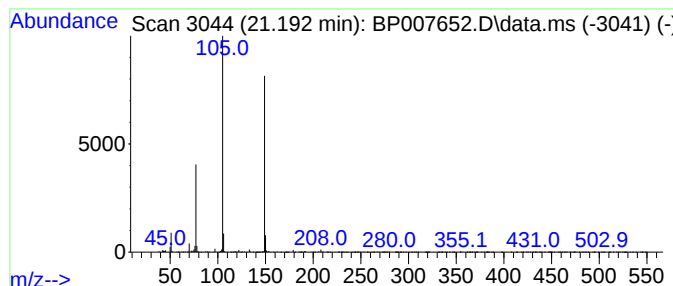
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3-Dioxolane, 2-(methoxyme... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.192	2.14 ng/ul	291368	Chrysene-d12	21.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	64		
2	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	64		
3	Phenyl(2-phenyl-1,3-dioxolan-2-y...	256	C16H16O3	005694-69-9	64		
4	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	64		
5	Benzoic acid, 2-(3-nitrophenyl)e...	271	C15H13NO4	1000368-22-4	64		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Dioxolane, ...	21.192	2.1	ng/ul	291368	5	21.416	2720710	20.0