

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102523\
 Data File : BP017822.D
 Acq On : 26 Oct 2023 02:38
 Operator : MA/JU
 Sample : 04927-16DL 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCD20DL

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-BP101823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017822.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.740	106	111	118	rBV3	13527	38393	1.44%	0.282%
2	4.399	219	223	227	rVB2	8970	14128	0.53%	0.104%
3	5.499	407	410	411	rBV3	2327	2702	0.10%	0.020%
4	5.517	411	413	419	rVB7	2202	4195	0.16%	0.031%
5	6.028	496	500	502	rVB5	2396	3223	0.12%	0.024%
6	6.058	502	505	508	rBV5	2678	3243	0.12%	0.024%
7	6.481	572	577	580	rVV7	4810	6980	0.26%	0.051%
8	6.628	598	602	603	rBV4	2502	3250	0.12%	0.024%
9	6.658	604	607	612	rVV6	4248	6932	0.26%	0.051%
10	6.846	635	639	644	rVB2	11237	20263	0.76%	0.149%
11	7.058	669	675	681	rBV2	24269	52833	1.99%	0.388%
12	7.116	681	685	693	rVB2	23352	42055	1.58%	0.308%
13	7.234	700	705	720	rVB	29804	61233	2.30%	0.449%
14	7.411	729	735	745	rBV	37444	73464	2.76%	0.539%
15	7.752	790	793	795	rBV4	2013	2708	0.10%	0.020%
16	7.881	808	815	826	rBV	587602	965273	36.28%	7.080%
17	8.616	932	940	953	rBV3	29482	84705	3.18%	0.621%
18	9.093	1015	1021	1028	rBV	28687	55461	2.08%	0.407%
19	9.269	1048	1051	1053	rBV4	5570	5885	0.22%	0.043%
20	9.322	1058	1060	1063	rVB4	3688	3646	0.14%	0.027%
21	9.810	1138	1143	1149	rBV3	22383	42925	1.61%	0.315%
22	9.952	1162	1167	1170	rVB6	1649	3058	0.11%	0.022%
23	10.352	1230	1235	1243	rBV2	23865	61754	2.32%	0.453%
24	10.710	1287	1296	1311	rBV	738088	1246647	46.86%	9.144%
25	10.922	1327	1332	1333	rBV5	2773	2999	0.11%	0.022%
26	11.704	1459	1465	1467	rBV7	1658	2662	0.10%	0.020%
27	11.746	1467	1472	1479	rBV8	4838	13988	0.53%	0.103%
28	12.069	1521	1527	1529	rBV7	2288	3776	0.14%	0.028%
29	12.457	1589	1593	1596	rBV6	2721	3254	0.12%	0.024%
30	12.757	1636	1644	1647	rBV8	7476	15163	0.57%	0.111%
31	12.981	1679	1682	1685	rVB4	2459	3201	0.12%	0.023%
32	13.093	1695	1701	1705	rBV8	5053	9655	0.36%	0.071%
33	13.140	1705	1709	1713	rVB6	8947	12453	0.47%	0.091%
34	13.340	1737	1743	1745	rVB7	1807	3805	0.14%	0.028%
35	13.410	1750	1755	1760	rVV4	14225	22758	0.86%	0.167%
36	13.522	1773	1774	1780	rVB6	3941	5058	0.19%	0.037%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Title : SVOA CALIBRATION

37	13.569	1780	1782	1785	rBV4	2028	2736	0.10%	0.020%
38	13.681	1797	1801	1806	rBV7	4988	8312	0.31%	0.061%
39	13.793	1817	1820	1823	rVB5	2927	3780	0.14%	0.028%
40	13.993	1847	1854	1860	rBV	77469	113114	4.25%	0.830%
41	14.046	1860	1863	1868	rVB7	3055	5399	0.20%	0.040%
42	14.275	1897	1902	1912	rBV2	79595	132295	4.97%	0.970%
43	14.434	1921	1929	1933	rBV10	6888	15612	0.59%	0.115%
44	14.575	1946	1953	1963	rBV	1007188	1462718	54.98%	10.729%
45	14.875	2000	2004	2010	rBV9	8062	18006	0.68%	0.132%
46	14.969	2016	2020	2025	rBV8	10906	18792	0.71%	0.138%
47	15.081	2036	2039	2043	rVB5	3937	4916	0.18%	0.036%
48	15.128	2043	2047	2052	rBV8	9047	15404	0.58%	0.113%
49	15.210	2052	2061	2068	rBV2	51217	105126	3.95%	0.771%
50	15.369	2084	2088	2092	rVB7	4638	7761	0.29%	0.057%
51	15.398	2092	2093	2095	rBV2	3084	2752	0.10%	0.020%
52	15.451	2097	2102	2103	rBV5	4003	4787	0.18%	0.035%
53	15.481	2105	2107	2112	rVB6	3430	3575	0.13%	0.026%
54	15.581	2118	2124	2133	rBV	121741	192818	7.25%	1.414%
55	15.740	2147	2151	2160	rVB9	15385	28171	1.06%	0.207%
56	16.398	2261	2263	2264	rBV2	5495	3400	0.13%	0.025%
57	16.992	2358	2364	2378	rBV	1863210	2660509	100.00%	19.515%
58	17.139	2384	2389	2397	rVB2	99326	162080	6.09%	1.189%
59	17.369	2422	2428	2438	rBV	1170458	1661521	62.45%	12.187%
60	17.469	2440	2445	2450	rBV2	117483	171929	6.46%	1.261%
61	18.222	2569	2573	2578	rBV2	72284	121173	4.55%	0.889%
62	19.728	2826	2829	2837	rVB	163579	226729	8.52%	1.663%
63	21.510	3128	3132	3138	rVB	1396585	1725815	64.87%	12.659%
64	21.645	3153	3155	3160	rVB	155433	174433	6.56%	1.279%
65	24.063	3559	3566	3574	rVB	817090	1676103	63.00%	12.294%

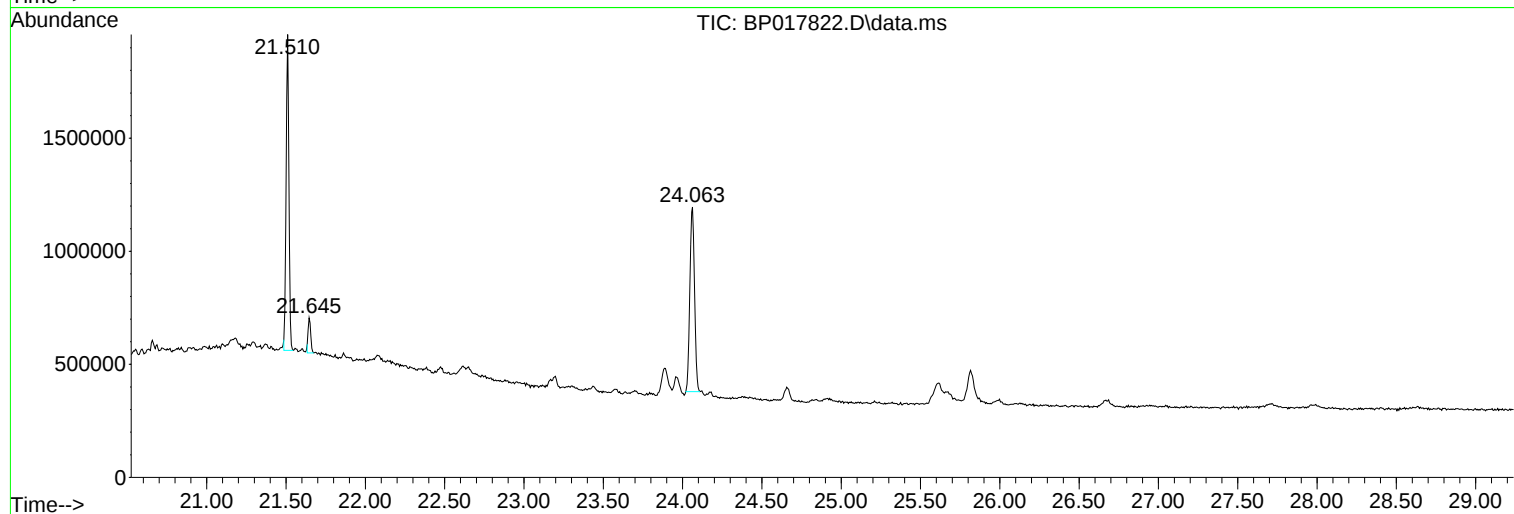
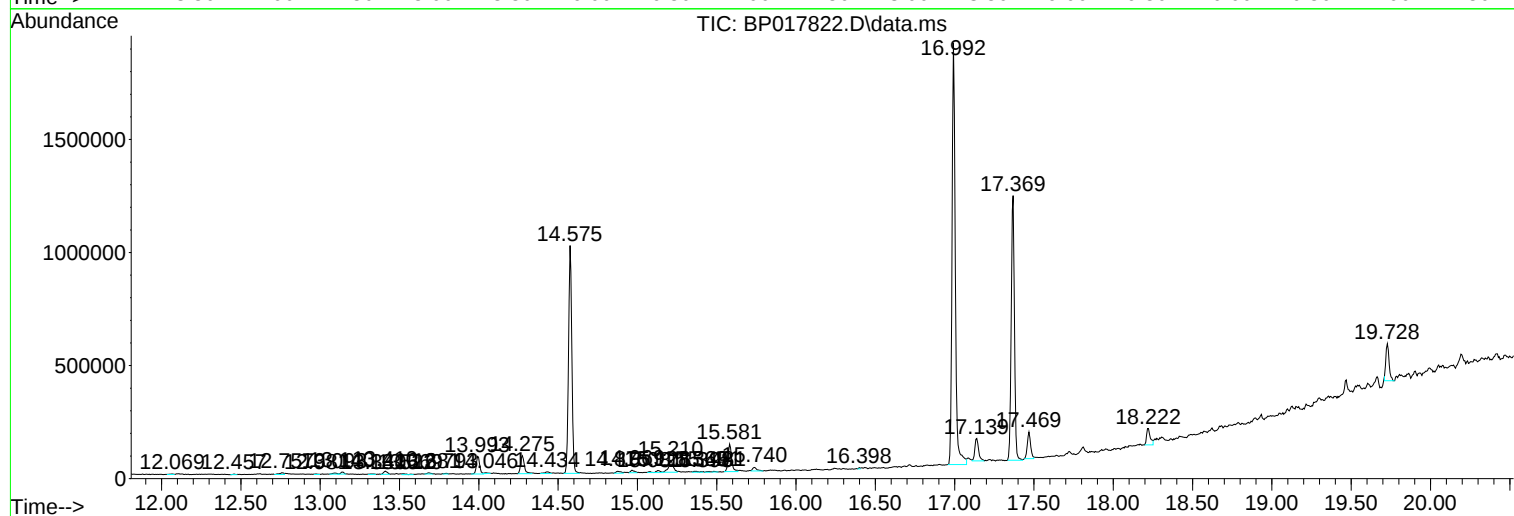
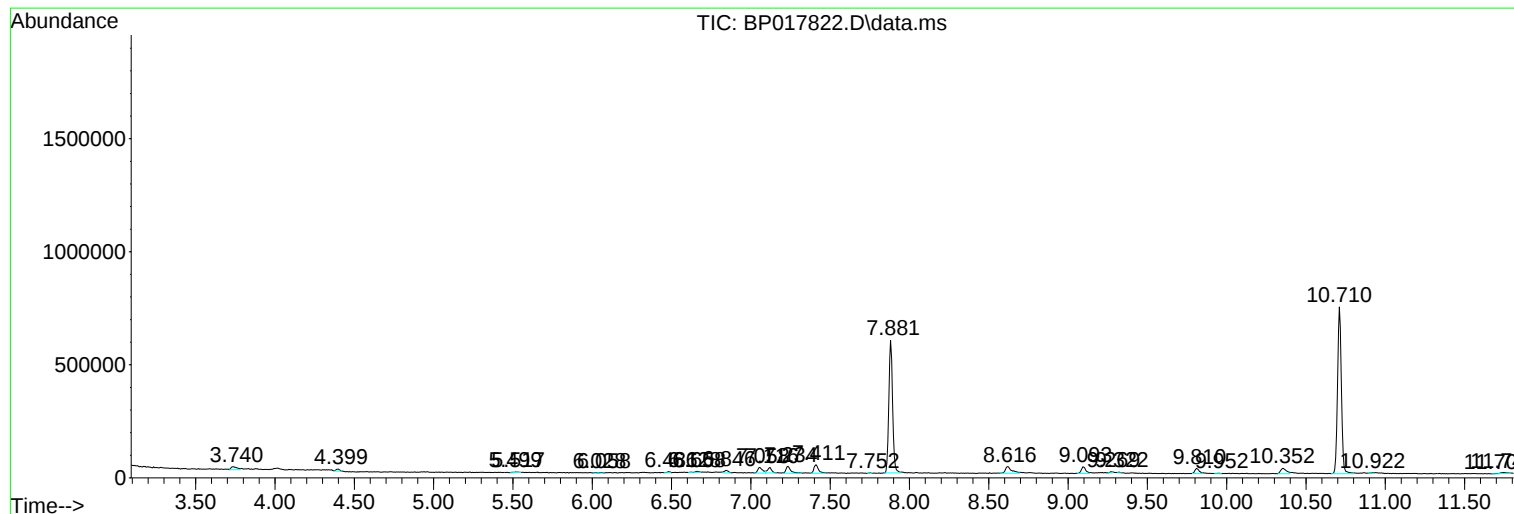
Sum of corrected areas: 13633494

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

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



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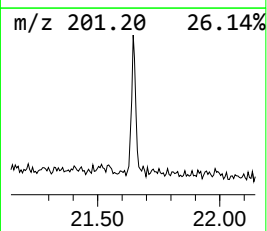
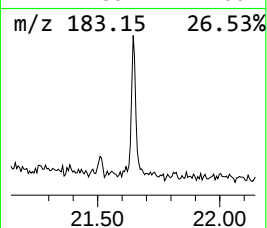
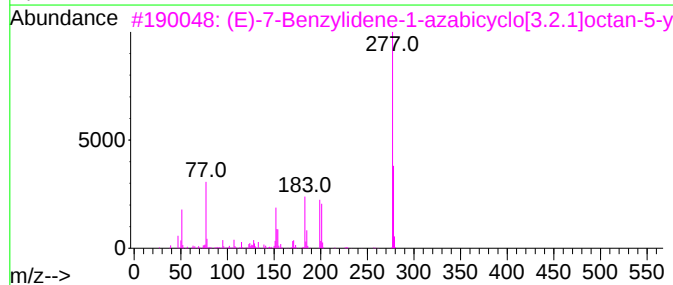
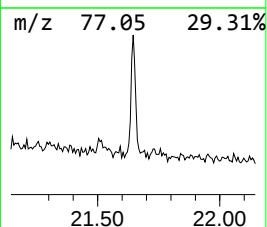
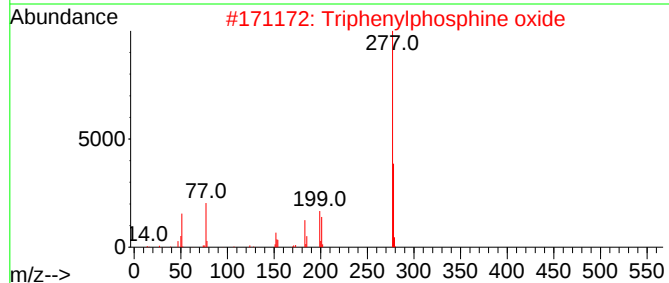
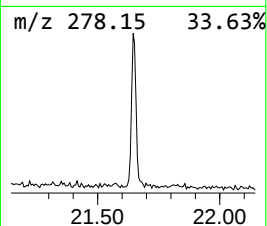
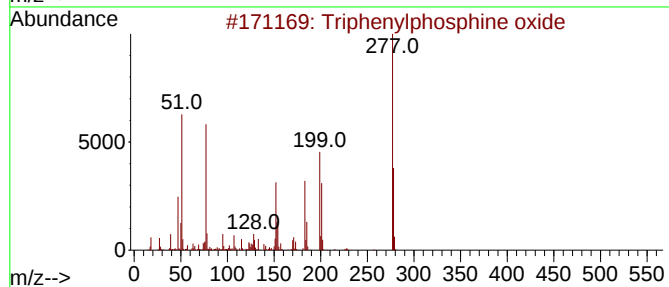
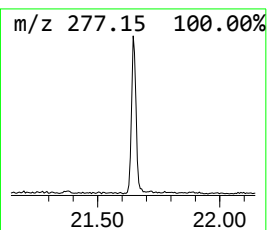
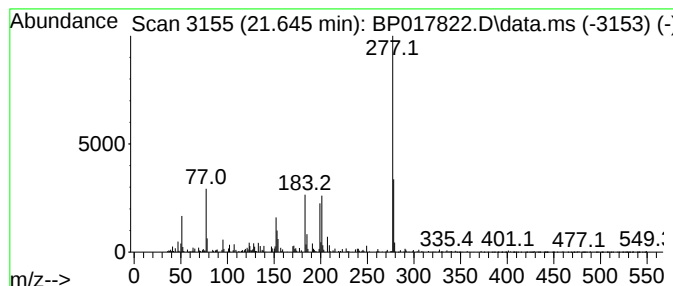
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.645	2.02 ng/ul	174433	Chrysene-d12	21.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	93
2			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	91
3			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	87
4			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	58
5			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	53



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Triphenylphosph...	21.645	2.0	ng/u1	174433	5	21.510	1725820	20.0