

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
 Data File : BP006568.D
 Acq On : 07 Aug 2021 14:58
 Operator : CG/JU
 Sample : M3163-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGEY4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: 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-BP072421.M

Title : SVOA CALIBRATION

Signal : TIC: BP006568.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.275	29	32	36	rVB2	37152	45159	1.48%	0.116%
2	3.681	97	101	113	rBV	324848	506556	16.61%	1.303%
3	6.928	647	653	665	rVB	608502	1001182	32.83%	2.575%
4	7.099	676	682	691	rVB	489746	799482	26.22%	2.056%
5	7.281	707	713	722	rVB	617614	998754	32.75%	2.568%
6	7.751	787	793	801	rVB	684800	1069460	35.07%	2.750%
7	8.457	907	913	921	rBV	635050	1007587	33.04%	2.591%
8	8.569	930	932	940	rVB3	12117	20140	0.66%	0.052%
9	8.904	983	989	1000	rVB	627026	1002169	32.86%	2.577%
10	9.022	1007	1009	1014	rVB6	5551	7820	0.26%	0.020%
11	9.622	1104	1111	1120	rBV	477041	778529	25.53%	2.002%
12	10.157	1195	1202	1214	rBV	647361	1065059	34.92%	2.739%
13	10.328	1225	1231	1243	rBV	121882	217023	7.12%	0.558%
14	10.534	1258	1266	1273	rBV	921289	1513822	49.64%	3.893%
15	10.675	1283	1290	1305	rBV	627705	1099123	36.04%	2.826%
16	11.016	1345	1348	1350	rBV4	2665	3209	0.11%	0.008%
17	11.604	1444	1448	1451	rBV6	2906	4706	0.15%	0.012%
18	11.810	1479	1483	1491	rVB4	8339	17691	0.58%	0.045%
19	11.963	1505	1509	1512	rVB5	2912	4851	0.16%	0.012%
20	12.045	1517	1523	1525	rBV7	3956	7133	0.23%	0.018%
21	12.122	1532	1536	1541	rVB2	12323	20820	0.68%	0.054%
22	12.322	1567	1570	1571	rBV2	2936	3223	0.11%	0.008%
23	12.716	1632	1637	1639	rBV6	3440	3430	0.11%	0.009%
24	12.739	1639	1641	1646	rVB5	2966	4903	0.16%	0.013%
25	12.992	1683	1684	1689	rVB5	4212	4223	0.14%	0.011%
26	13.104	1698	1703	1707	rVB8	1888	3530	0.12%	0.009%
27	13.216	1717	1722	1727	rVB7	7588	13258	0.43%	0.034%
28	13.292	1730	1735	1737	rVV6	4157	6392	0.21%	0.016%
29	13.345	1741	1744	1749	rVV7	4659	8738	0.29%	0.022%
30	13.392	1749	1752	1758	rVB8	2635	5159	0.17%	0.013%
31	13.798	1813	1821	1832	rBV	1255416	1753975	57.52%	4.511%
32	14.069	1857	1867	1880	rBV	1357940	1960314	64.28%	5.041%
33	14.298	1902	1906	1908	rVB5	3724	3917	0.13%	0.010%
34	14.381	1912	1920	1936	rBV	1306096	1896121	62.18%	4.876%
35	14.586	1949	1955	1976	rBV	386519	705335	23.13%	1.814%
36	14.745	1980	1982	1989	rVV7	6615	9602	0.31%	0.025%

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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-BP072421.M
 Title : SVOA CALIBRATION

37	14.810	1989	1993	1998	rVB7	8834	12047	0.40%	0.031%
38	15.198	2054	2059	2060	rBV5	3785	5418	0.18%	0.014%
39	15.298	2073	2076	2081	rBV6	2812	5240	0.17%	0.013%
40	15.375	2082	2089	2103	rVV	1746162	2476411	81.21%	6.368%
41	15.492	2104	2109	2125	rVV	406179	594688	19.50%	1.529%
42	15.616	2126	2130	2134	rVB3	19030	27257	0.89%	0.070%
43	15.833	2163	2167	2170	rVB6	5422	7024	0.23%	0.018%
44	16.063	2200	2206	2212	rBV6	4695	8392	0.28%	0.022%
45	16.169	2220	2224	2227	rVB5	3669	5694	0.19%	0.015%
46	16.551	2285	2289	2292	rVB5	2849	4453	0.15%	0.011%
47	16.686	2305	2312	2319	rVV5	9529	24436	0.80%	0.063%
48	16.810	2329	2333	2336	rVB6	4407	5997	0.20%	0.015%
49	16.916	2347	2351	2355	rBV4	6788	11589	0.38%	0.030%
50	16.945	2355	2356	2363	rVB6	4794	6127	0.20%	0.016%
51	17.004	2363	2366	2368	rBV4	3353	3646	0.12%	0.009%
52	17.075	2376	2378	2380	rBV3	3039	3479	0.11%	0.009%
53	17.133	2381	2388	2398	rVV	1528557	2175025	71.32%	5.593%
54	17.227	2398	2404	2416	rVV2	1802621	2614037	85.72%	6.722%
55	17.822	2500	2505	2508	rBV3	17038	23096	0.76%	0.059%
56	17.910	2519	2520	2524	rBV4	2539	3708	0.12%	0.010%
57	18.033	2536	2541	2550	rBV	123348	202821	6.65%	0.522%
58	18.298	2584	2586	2588	rBV3	3545	3301	0.11%	0.008%
59	18.333	2588	2592	2595	rVV6	5005	8838	0.29%	0.023%
60	18.451	2607	2612	2617	rBV	30995	48930	1.60%	0.126%
61	18.898	2685	2688	2693	rBV7	12017	20558	0.67%	0.053%
62	19.051	2710	2714	2718	rBV2	27518	33043	1.08%	0.085%
63	19.116	2722	2725	2728	rBV	23467	31340	1.03%	0.081%
64	19.274	2748	2752	2760	rBV3	59764	87562	2.87%	0.225%
65	19.474	2780	2786	2793	rBV	2300579	3049572	100.00%	7.842%
66	20.463	2950	2954	2959	rBV	1161202	1218634	39.96%	3.134%
67	20.957	3035	3038	3046	rBV3	124393	202070	6.63%	0.520%
68	21.227	3079	3084	3095	rBV2	1981576	2493126	81.75%	6.411%
69	22.868	3359	3363	3367	rVB	141400	208912	6.85%	0.537%
70	23.404	3447	3454	3462	rBV2	1602708	2983566	97.84%	7.673%
71	23.557	3473	3480	3489	rVB	1415013	2704005	88.67%	6.954%

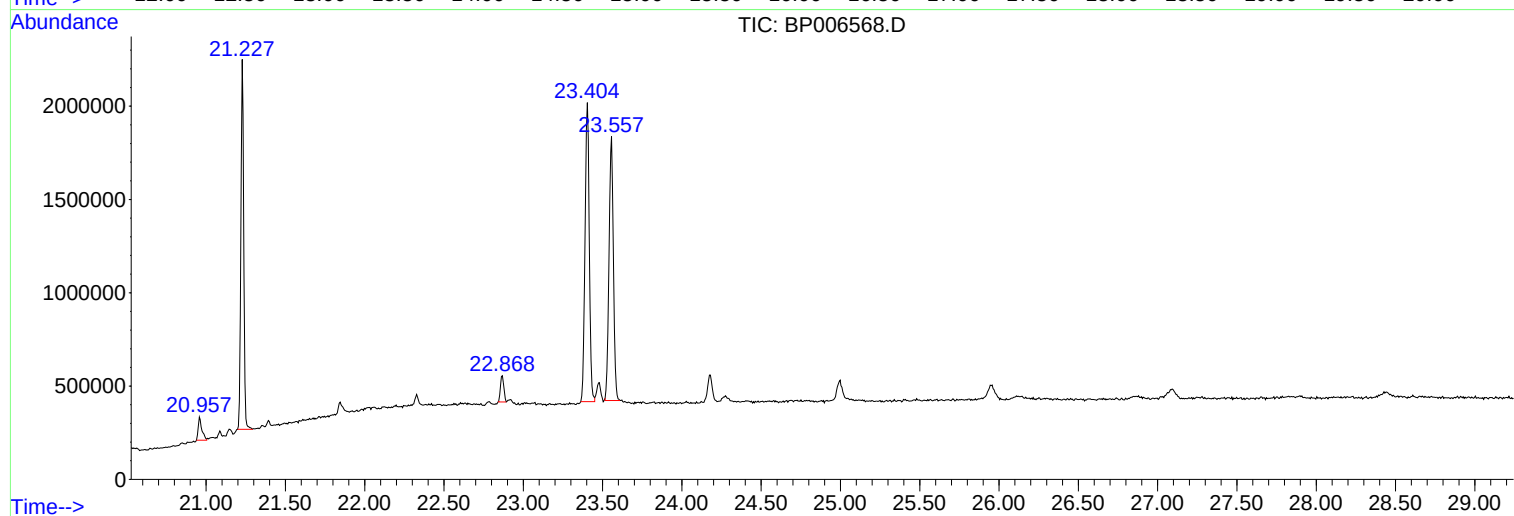
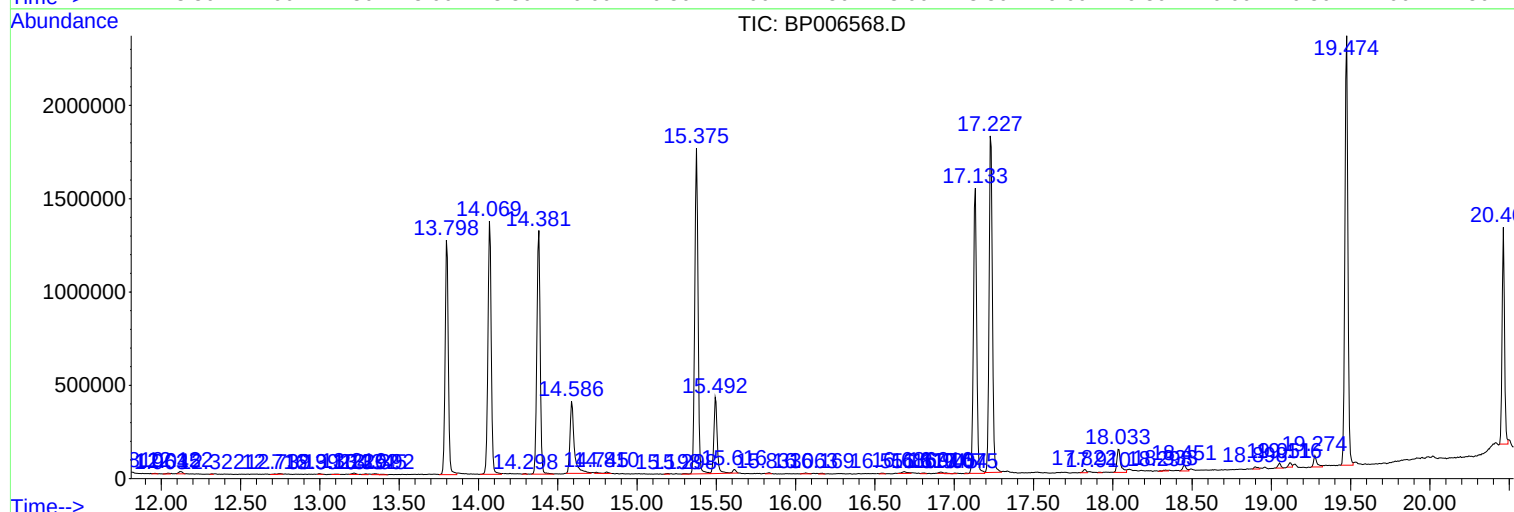
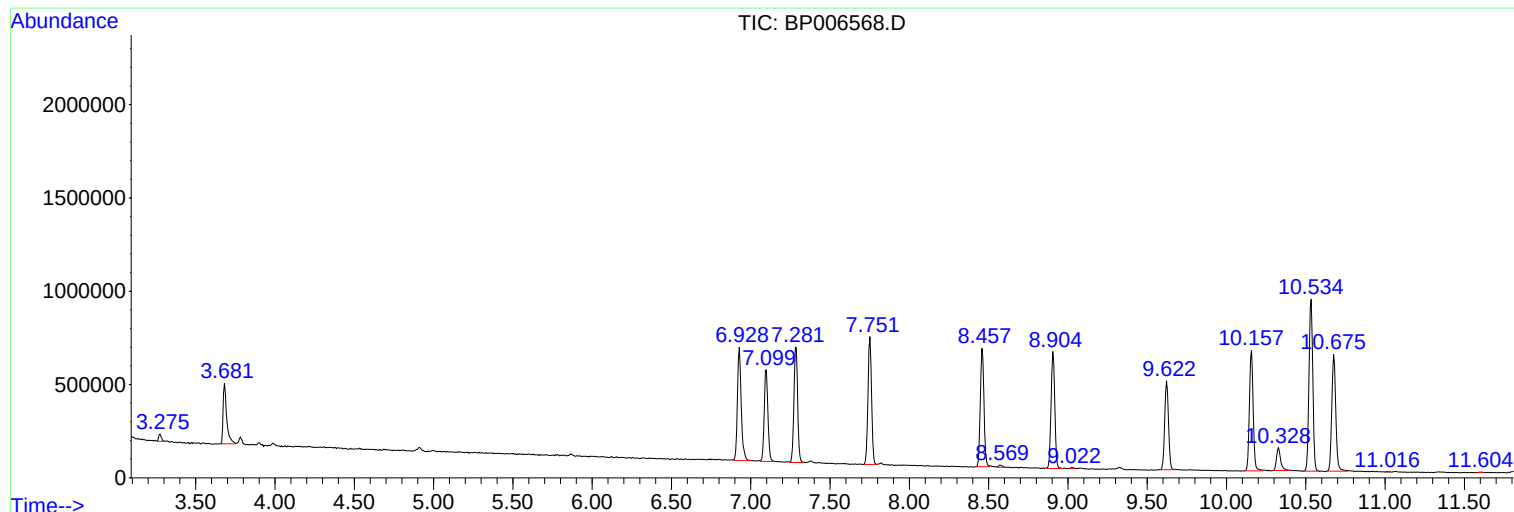
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ClientSampleId :
BGEY4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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Sample : M3163-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGEY4

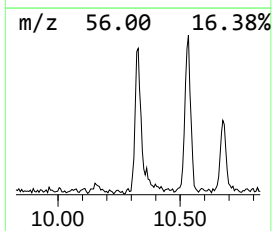
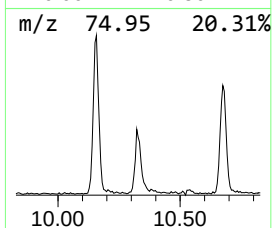
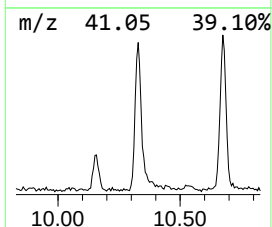
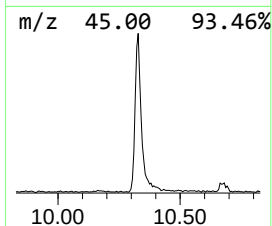
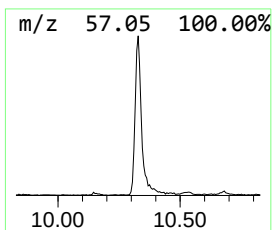
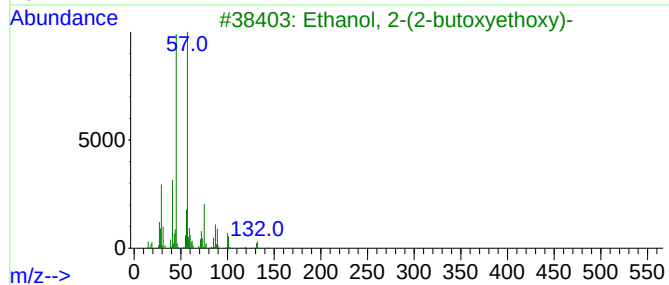
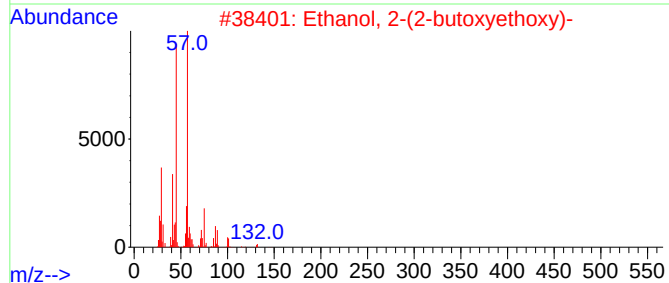
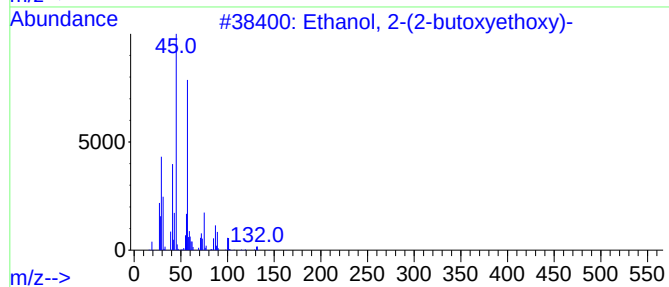
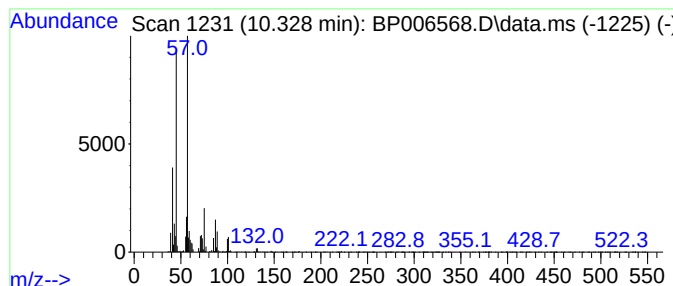
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.330	2.87 ng/ul	217023	Naphthalene-d8	10.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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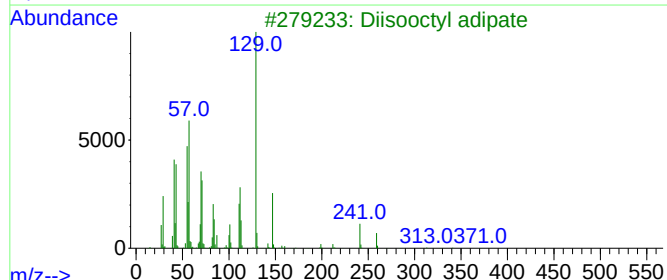
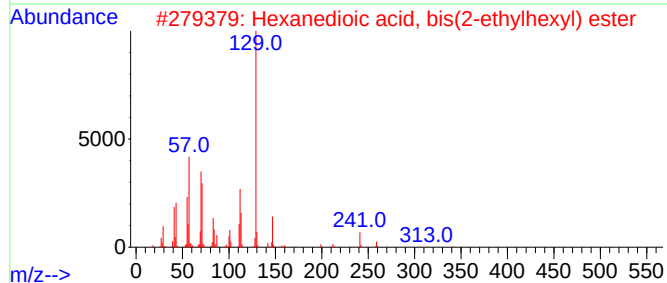
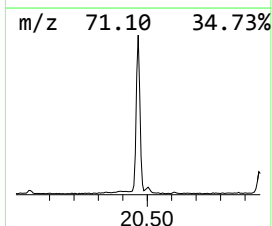
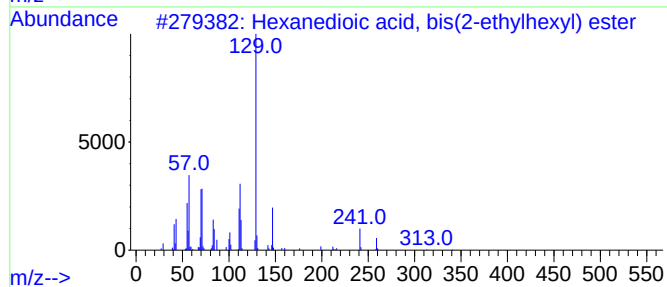
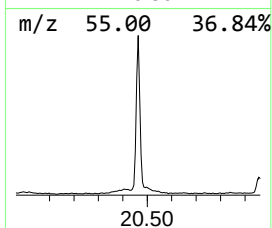
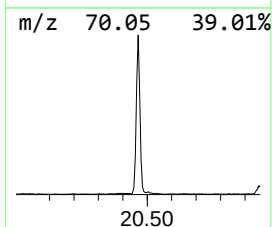
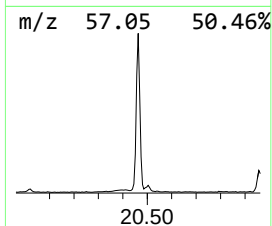
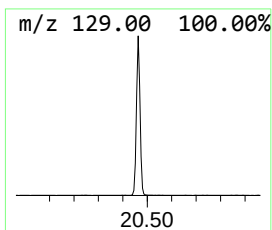
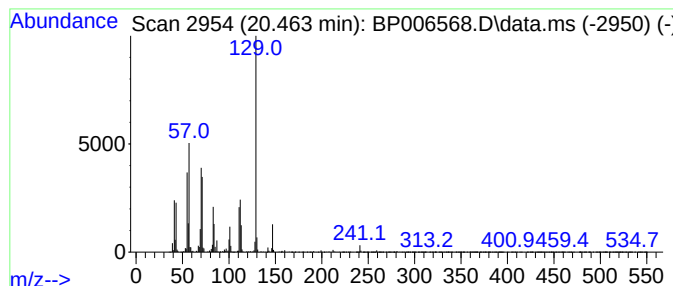
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexanedioic acid, bis(2-eth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.460	9.78 ng/ul	1218630	Chrysene-d12	21.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
3			Diisooctyl adipate	370	C22H42O4	001330-86-5	90
4			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
5			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	86



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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
Ethanol, 2-(2-b...	10.330	2.9	ng/ul	217023	2	10.534	1513820	20.0
Hexanedioic aci...	20.460	9.8	ng/ul	1218630	5	21.227	2493130	20.0