

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
 Data File : BP021183.D
 Acq On : 27 Jul 2024 00:40
 Operator : MA/JU
 Sample : P3300-23
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1GD2

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021183.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.175	29	32	41	rVB	33982	49079	1.15%	0.124%
2	3.605	103	105	122	rBV	127837	240573	5.63%	0.606%
3	3.899	151	155	161	rVB	21380	33916	0.79%	0.085%
4	4.275	214	219	221	rVB5	3032	4557	0.11%	0.011%
5	4.799	306	308	312	rVB3	4701	4567	0.11%	0.011%
6	5.234	377	382	387	rVB6	4570	9616	0.22%	0.024%
7	5.740	463	468	472	rBV6	5819	10561	0.25%	0.027%
8	5.934	496	501	508	rVB6	4127	8344	0.20%	0.021%
9	6.328	560	568	573	rBV10	3801	8441	0.20%	0.021%
10	6.752	635	640	642	rBV4	3726	5708	0.13%	0.014%
11	6.875	652	661	676	rBV	141490	321596	7.52%	0.810%
12	6.993	676	681	694	rVB	529856	822710	19.24%	2.071%
13	7.199	709	716	731	rBV	566912	1044713	24.43%	2.630%
14	7.646	781	792	799	rBV	615532	985103	23.04%	2.480%
15	7.710	799	803	815	rVB2	41321	101687	2.38%	0.256%
16	7.934	836	841	847	rBV8	7000	12589	0.29%	0.032%
17	8.381	909	917	929	rBV	469641	866138	20.26%	2.180%
18	8.475	929	933	941	rVB	42015	79662	1.86%	0.201%
19	8.734	970	977	982	rBV2	35251	62258	1.46%	0.157%
20	8.799	982	988	1009	rVV	561541	1057359	24.73%	2.662%
21	8.957	1011	1015	1023	rVB8	6756	13790	0.32%	0.035%
22	9.510	1103	1109	1125	rBV	479010	885100	20.70%	2.228%
23	9.834	1159	1164	1168	rBV8	3574	6998	0.16%	0.018%
24	9.887	1168	1173	1177	rVV8	5237	10426	0.24%	0.026%
25	10.075	1195	1205	1221	rBV	588709	1298199	30.36%	3.268%
26	10.263	1233	1237	1242	rVB7	8448	14285	0.33%	0.036%
27	10.334	1242	1249	1255	rBV4	18604	37769	0.88%	0.095%
28	10.416	1255	1263	1270	rVV	756306	1304083	30.50%	3.283%
29	10.475	1270	1273	1282	rVB7	20227	36965	0.86%	0.093%
30	10.569	1282	1289	1309	rBV	543400	1098377	25.69%	2.765%
31	10.951	1350	1354	1359	rVB8	3403	5101	0.12%	0.013%
32	12.028	1528	1537	1543	rBV3	8508	22525	0.53%	0.057%
33	12.616	1631	1637	1640	rBV7	4263	8571	0.20%	0.022%
34	12.798	1662	1668	1671	rVV8	4594	8529	0.20%	0.021%
35	12.863	1673	1679	1688	rVB7	13187	26940	0.63%	0.068%
36	13.128	1720	1724	1727	rBV3	7377	10774	0.25%	0.027%

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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-BP071024.MA.M
 Title : SVOA CALIBRATION

37	13.245	1740	1744	1746	rBV5	4090	5581	0.13%	0.014%
38	13.504	1779	1788	1793	rBV5	4982	8887	0.21%	0.022%
39	13.692	1814	1820	1834	rBV	1330657	1904139	44.53%	4.793%
40	13.787	1834	1836	1840	rVB5	4883	6430	0.15%	0.016%
41	13.975	1860	1868	1883	rBV	1443918	2259051	52.83%	5.687%
42	14.204	1904	1907	1910	rBV5	4721	6901	0.16%	0.017%
43	14.287	1913	1921	1929	rVV	1017345	1732521	40.52%	4.361%
44	14.369	1929	1935	1940	rVB4	21475	37076	0.87%	0.093%
45	14.622	1973	1978	1983	rBV2	44930	100298	2.35%	0.252%
46	14.769	2001	2003	2008	rBV6	3419	4454	0.10%	0.011%
47	14.928	2026	2030	2033	rBV5	3851	6306	0.15%	0.016%
48	14.986	2034	2040	2048	rVB4	11595	27824	0.65%	0.070%
49	15.081	2053	2056	2060	rVB3	6974	10243	0.24%	0.026%
50	15.134	2060	2065	2070	rBV3	8112	12804	0.30%	0.032%
51	15.298	2086	2093	2110	rVV	1895556	2843148	66.50%	7.157%
52	15.475	2115	2123	2132	rBV	377616	842968	19.72%	2.122%
53	15.681	2154	2158	2162	rBV	19835	24389	0.57%	0.061%
54	15.886	2191	2193	2197	rVB5	5889	7032	0.16%	0.018%
55	16.675	2325	2327	2332	rVB5	5720	8052	0.19%	0.020%
56	17.057	2389	2392	2394	rBV4	5887	7250	0.17%	0.018%
57	17.104	2394	2400	2410	rVV	1323278	2045192	47.83%	5.148%
58	17.204	2410	2417	2435	rVB	2123169	3186575	74.53%	8.022%
59	17.398	2447	2450	2457	rVB	10951	14265	0.33%	0.036%
60	17.786	2510	2516	2522	rBV	330516	455272	10.65%	1.146%
61	18.027	2553	2557	2567	rBV5	32186	76337	1.79%	0.192%
62	18.104	2567	2570	2579	rVB	21112	33719	0.79%	0.085%
63	19.133	2741	2745	2749	rVB	33785	41999	0.98%	0.106%
64	19.369	2782	2785	2791	rBV6	29145	53309	1.25%	0.134%
65	19.586	2816	2822	2839	rBV2	2435222	3951903	92.43%	9.948%
66	21.551	3150	3156	3166	rBV2	1332556	2392522	55.96%	6.023%
67	24.633	3672	3680	3692	rVB	1427841	4275699	100.00%	10.763%
68	24.868	3712	3720	3738	rVB	987287	2825400	66.08%	7.112%

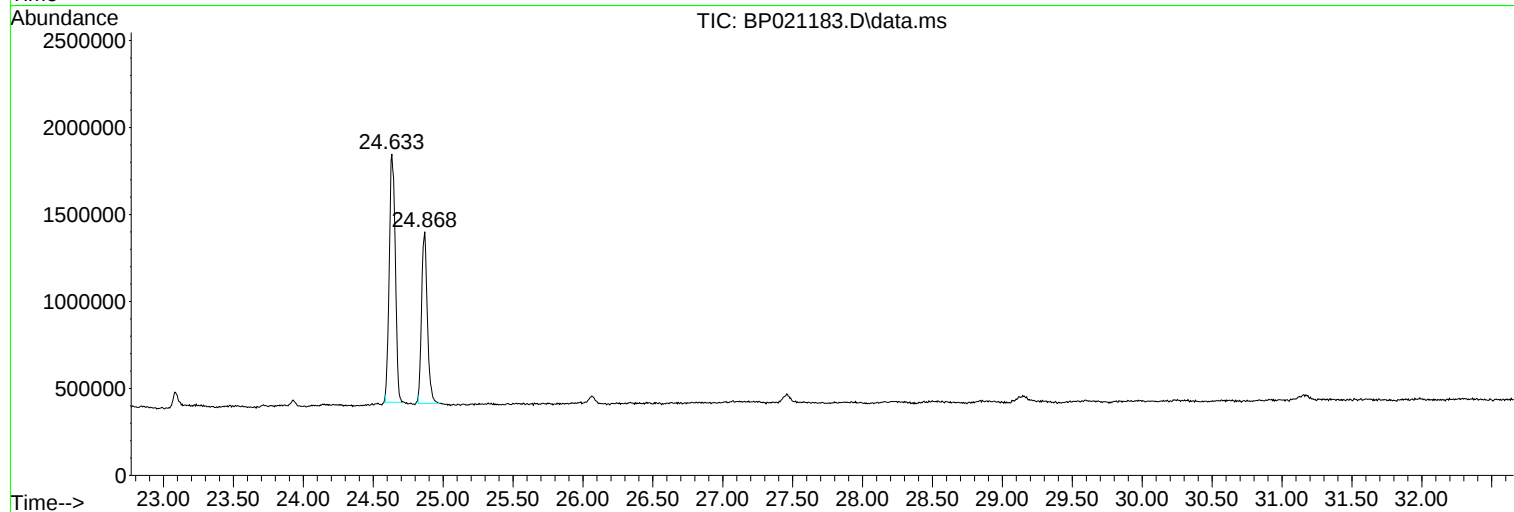
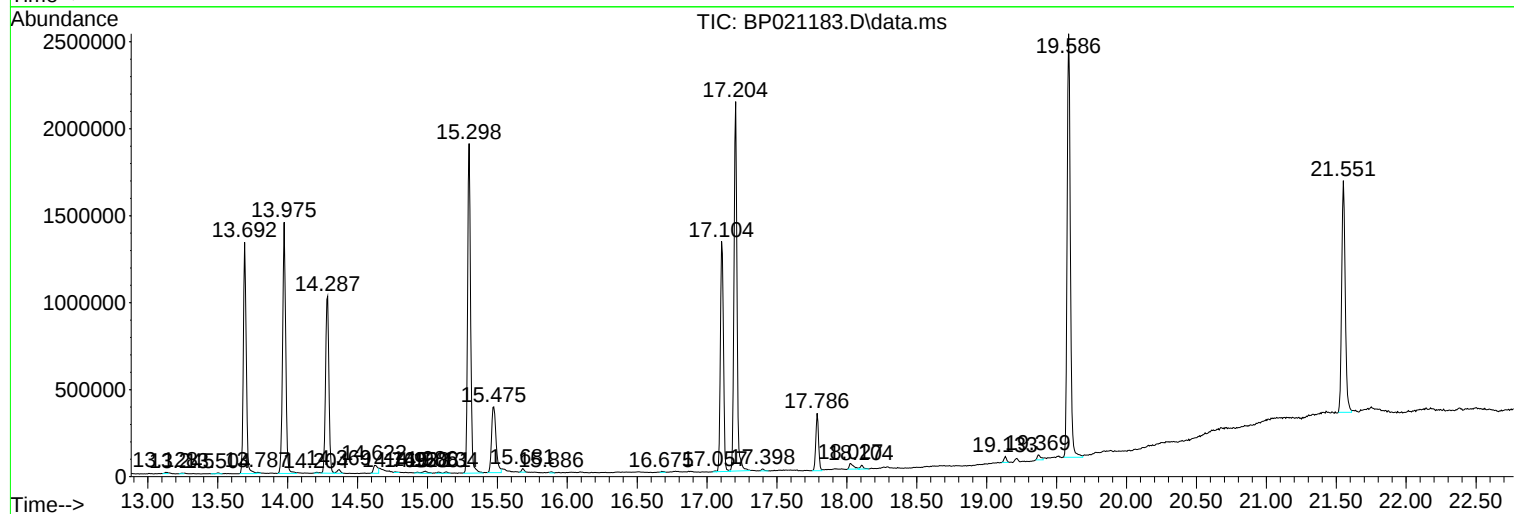
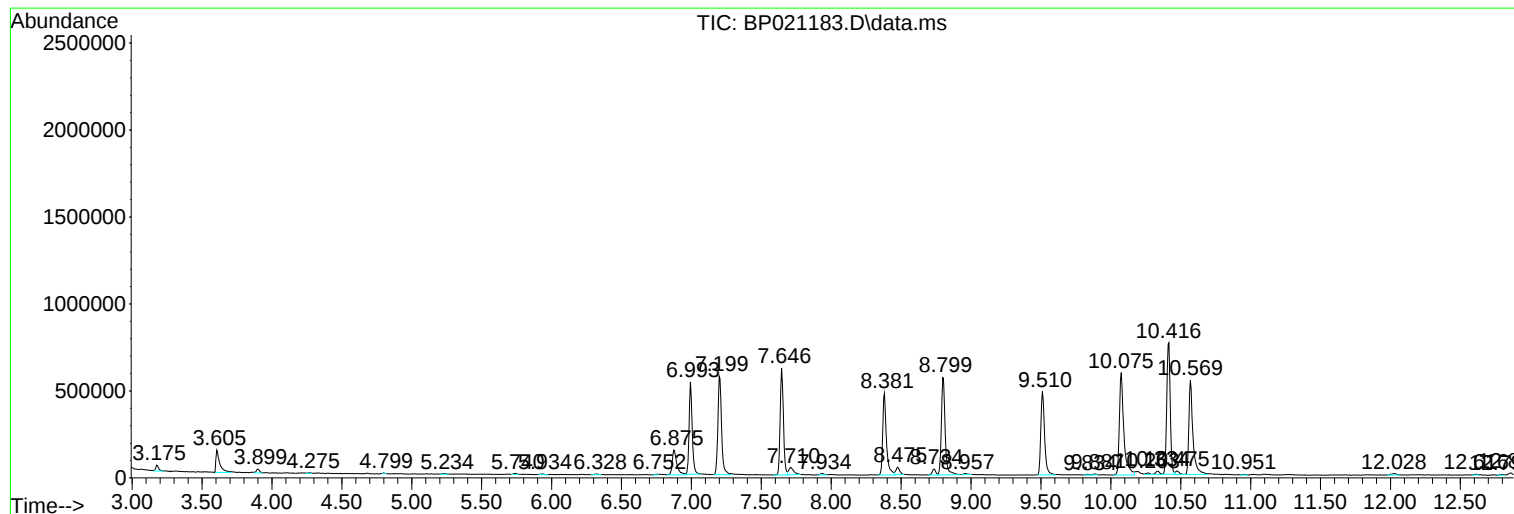
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : MA/JU
Sample : P3300-23
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1GD2

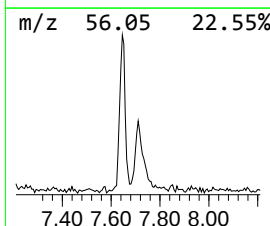
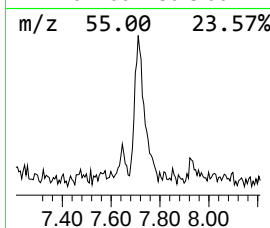
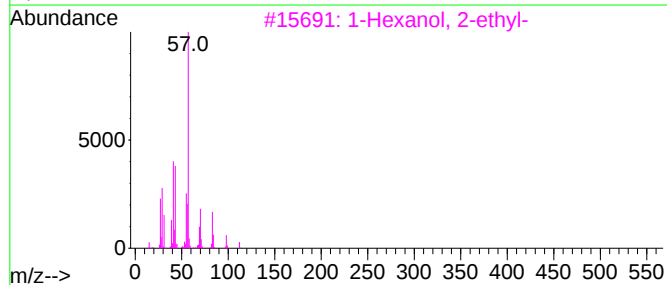
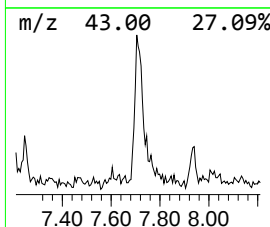
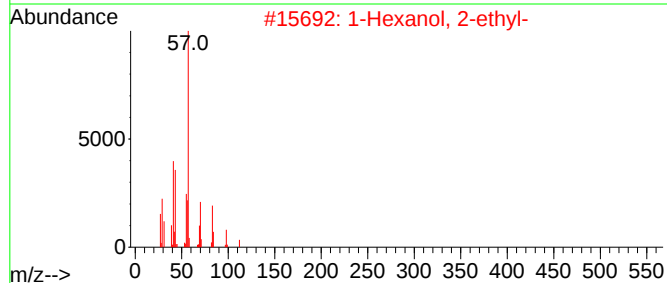
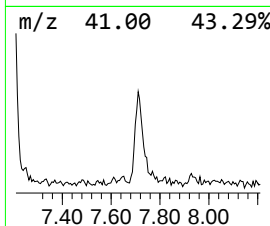
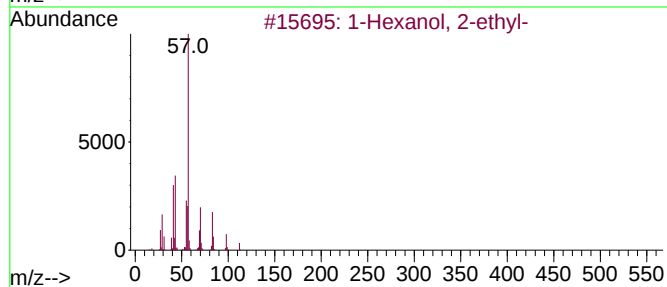
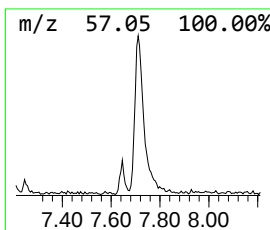
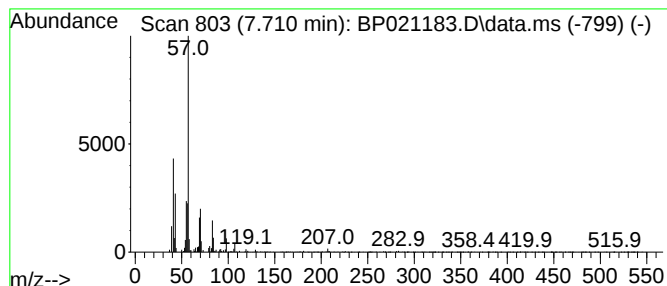
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.710	2.06 ng/ul	101687	1,4-Dichlorobenzene-d4	7.646

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	59



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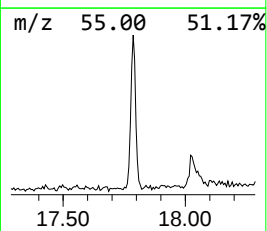
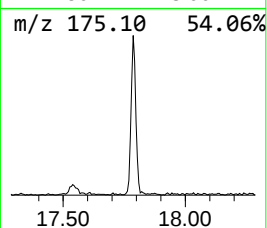
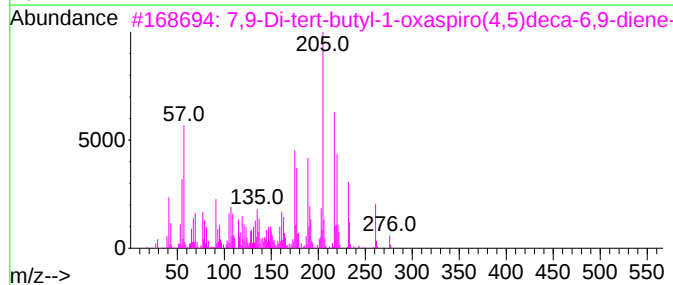
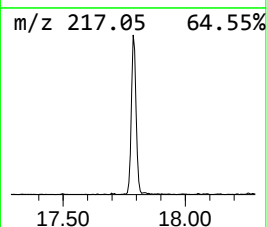
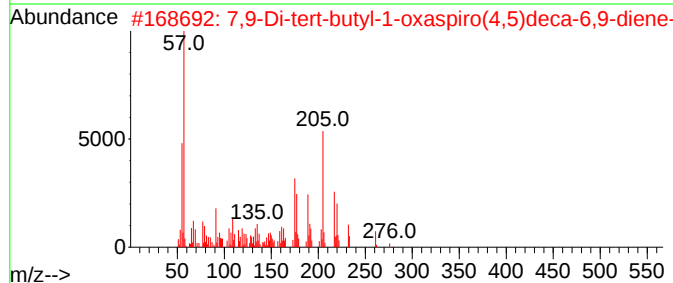
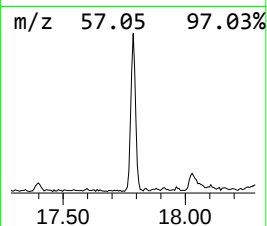
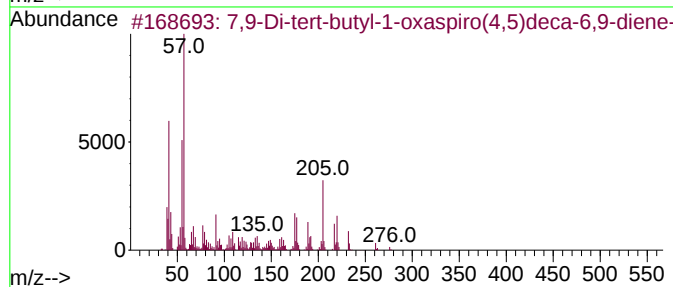
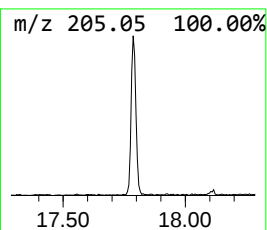
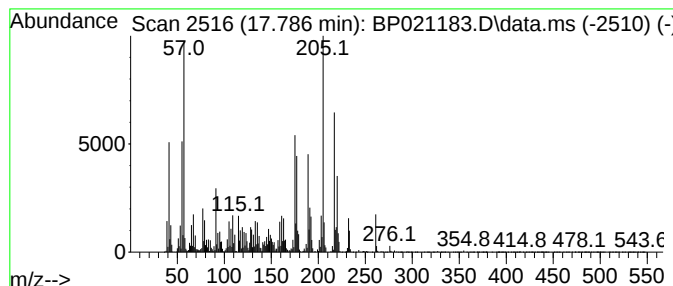
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.786	4.45 ng/ul	455272	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	98		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	91		
4	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	50		
5	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	50		



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	7.710	2.1	ng/ul	101687	1	7.646	985103	20.0
7,9-Di-tert-but...	17.786	4.5	ng/ul	455272	4	17.104	2045190	20.0