

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035981.D
 Acq On : 28 Jul 2022 12:17
 Operator : CG/JU
 Sample : N3874-08
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GBJ34

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_M\Methods\SFAM-EPA-BM072522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM035981.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.269	27	31	37	rVB	112707	145412	1.94%	0.059%
2	3.616	85	90	97	rVB2	54565	88588	1.18%	0.036%
3	3.687	99	102	118	rVB	1328940	1872341	25.02%	0.755%
4	4.022	155	159	165	rBV	39968	49947	0.67%	0.020%
5	4.552	247	249	260	rVB9	22248	44568	0.60%	0.018%
6	4.969	314	320	326	rVB	286981	402353	5.38%	0.162%
7	6.057	500	505	511	rBV	43202	66055	0.88%	0.027%
8	7.040	665	672	675	rBV	1672735	2924810	39.08%	1.180%
9	7.204	694	700	708	rBV	1393201	2121517	28.34%	0.856%
10	7.299	709	716	727	rVV	2604344	3779214	50.49%	1.525%
11	7.404	727	734	742	rVB	1658292	2674147	35.73%	1.079%
12	7.869	805	813	821	rBV	1287483	2021724	27.01%	0.816%
13	8.322	883	890	899	rBV	3879227	5809167	77.61%	2.343%
14	8.593	928	936	941	rBV	1935174	3153860	42.14%	1.272%
15	8.651	941	946	950	rVV	3850525	6050001	80.83%	2.441%
16	8.698	950	954	970	rVB	2572466	4197384	56.08%	1.693%
17	8.951	990	997	1004	rVB	1925868	3067931	40.99%	1.238%
18	9.040	1004	1012	1022	rBV	1550163	2456910	32.83%	0.991%
19	9.610	1102	1109	1120	rBV	2445024	3971742	53.07%	1.602%
20	9.763	1127	1135	1143	rBV	1235913	1969906	26.32%	0.795%
21	9.857	1144	1151	1170	rVB	2903623	4616225	61.68%	1.862%
22	10.098	1183	1192	1199	rBV	2533124	3826174	51.12%	1.543%
23	10.304	1219	1227	1242	rBV2	1866889	4657944	62.23%	1.879%
24	10.681	1280	1291	1295	rBV	1703016	2963421	39.59%	1.195%
25	10.728	1295	1299	1306	rVB	2168958	3392671	45.33%	1.369%
26	10.822	1307	1315	1326	rBV	1429716	2449950	32.73%	0.988%
27	10.957	1334	1338	1340	rBV	38106	55421	0.74%	0.022%
28	11.010	1340	1347	1366	rVB	3929129	6189599	82.70%	2.497%
29	11.386	1404	1411	1417	rBV	60291	108435	1.45%	0.044%
30	11.451	1417	1422	1429	rVB	98994	148861	1.99%	0.060%
31	11.622	1441	1451	1473	rBV	878210	2435465	32.54%	0.982%
32	12.204	1541	1550	1557	rBV	868916	1335667	17.85%	0.539%
33	12.269	1557	1561	1567	rVB2	41198	66092	0.88%	0.027%
34	12.345	1567	1574	1584	rVB2	59869	102465	1.37%	0.041%
35	12.469	1589	1595	1603	rBV	164976	274301	3.66%	0.111%
36	12.563	1603	1611	1618	rBV	4592435	7274796	97.20%	2.935%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Title : SVOA CALIBRATION

37	12.704	1624	1635	1643	rVB2	3166703	7195660	96.14%	2.903%
38	12.798	1645	1651	1653	rBV4	22349	46004	0.61%	0.019%
39	13.022	1681	1689	1705	rBV	2574537	3953338	52.82%	1.595%
40	13.392	1741	1752	1759	rBV	1943568	2914556	38.94%	1.176%
41	13.769	1809	1816	1817	rBV5	28923	48589	0.65%	0.020%
42	13.928	1838	1843	1847	rVV	2921803	3888059	51.95%	1.568%
43	13.975	1847	1851	1859	rVB	1834825	2388684	31.91%	0.964%
44	14.210	1881	1891	1893	rVV	3563014	5393838	72.07%	2.176%
45	14.233	1893	1895	1902	rVB	2353453	3058424	40.86%	1.234%
46	14.427	1916	1928	1935	rBV	2508257	3642464	48.67%	1.469%
47	14.510	1936	1942	1948	rVV	2351280	3556468	47.52%	1.435%
48	14.575	1948	1953	1958	rVV	4142784	6047373	80.80%	2.440%
49	14.627	1958	1962	1971	rVV	1265719	1856956	24.81%	0.749%
50	14.716	1971	1977	1988	rVV	1351636	1944336	25.98%	0.784%
51	14.810	1988	1993	2000	rVV	886850	1235394	16.51%	0.498%
52	14.910	2004	2010	2018	rVB	2148930	2969757	39.68%	1.198%
53	15.298	2071	2076	2084	rBV2	56850	133066	1.78%	0.054%
54	15.374	2085	2089	2093	rVB	36052	51277	0.69%	0.021%
55	15.504	2100	2111	2115	rBV	4412945	6563044	87.69%	2.648%
56	15.557	2115	2120	2123	rVV	4042069	6051090	80.85%	2.441%
57	15.592	2123	2126	2129	rVV	2664261	3580022	47.83%	1.444%
58	15.639	2129	2134	2146	rVV2	3276255	5481696	73.24%	2.211%
59	15.745	2147	2152	2157	rVB	54601	93983	1.26%	0.038%
60	15.921	2175	2182	2188	rVB4	36466	68972	0.92%	0.028%
61	16.216	2228	2232	2240	rVV	51106	101677	1.36%	0.041%
62	16.380	2255	2260	2265	rBV4	20567	41312	0.55%	0.017%
63	16.557	2284	2290	2294	rVV2	78872	137488	1.84%	0.055%
64	16.733	2314	2320	2327	rVV	263500	356915	4.77%	0.144%
65	16.904	2342	2349	2364	rBV	5288748	7285926	97.35%	2.939%
66	17.010	2364	2367	2370	rBV	90311	110225	1.47%	0.044%
67	17.098	2379	2382	2388	rVB3	33119	57722	0.77%	0.023%
68	17.257	2402	2409	2416	rBV	2537734	3564299	47.62%	1.438%
69	17.351	2416	2425	2428	rVV2	4042280	5913123	79.00%	2.385%
70	17.386	2428	2431	2438	rVB	4127148	5511549	73.64%	2.223%
71	17.663	2471	2478	2490	rBV	4287595	6000884	80.18%	2.421%
72	17.751	2490	2493	2497	rVV3	24742	43556	0.58%	0.018%
73	17.804	2498	2502	2508	rVV2	224823	333144	4.45%	0.134%
74	17.868	2509	2513	2518	rVB	175131	222779	2.98%	0.090%
75	18.151	2556	2561	2566	rBV6	27990	52588	0.70%	0.021%
76	18.210	2566	2571	2577	rVB2	130653	215029	2.87%	0.087%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	18.321	2587	2590	2599	rVB2	46819	73408	0.98%	0.030%
78	18.433	2605	2609	2613	rBV	121397	162148	2.17%	0.065%
79	18.480	2613	2617	2623	rVB	211734	267746	3.58%	0.108%
80	18.674	2645	2650	2655	rVV	153935	223996	2.99%	0.090%
81	18.751	2658	2663	2671	rVB	179474	257515	3.44%	0.104%
82	19.210	2738	2741	2745	rVB	50935	57612	0.77%	0.023%
83	19.310	2746	2758	2767	rBV	2703670	3663776	48.95%	1.478%
84	19.551	2794	2799	2807	rVB	156809	221814	2.96%	0.089%
85	19.645	2809	2815	2820	rBV	4809638	6163499	82.35%	2.486%
86	19.692	2820	2823	2828	rVB	137029	170444	2.28%	0.069%
87	20.245	2913	2917	2924	rBV	169369	213452	2.85%	0.086%
88	20.574	2969	2973	2980	rBV	2671328	3119647	41.68%	1.258%
89	20.645	2981	2985	2990	rVV	694402	787530	10.52%	0.318%
90	20.874	3021	3024	3029	rVB	182462	207555	2.77%	0.084%
91	21.356	3101	3106	3111	rBV	3425907	3787744	50.61%	1.528%
92	21.439	3115	3120	3123	rVV	2620635	3723383	49.75%	1.502%
93	21.474	3123	3126	3135	rVB	5114903	5859909	78.29%	2.364%
94	22.009	3214	3217	3222	rBV2	208420	256205	3.42%	0.103%
95	22.233	3250	3255	3260	rBV	960464	1727646	23.08%	0.697%
96	23.133	3402	3408	3420	rVB	4380570	7433510	99.32%	2.999%
97	23.656	3490	3497	3501	rBV2	3009469	6007157	80.26%	2.423%
98	23.709	3501	3506	3515	rVV	3918626	7484626	100.00%	3.019%
99	23.809	3517	3523	3534	rVB2	1778443	3474972	46.43%	1.402%
100	26.268	3934	3941	3954	rVB	1214031	3671766	49.06%	1.481%

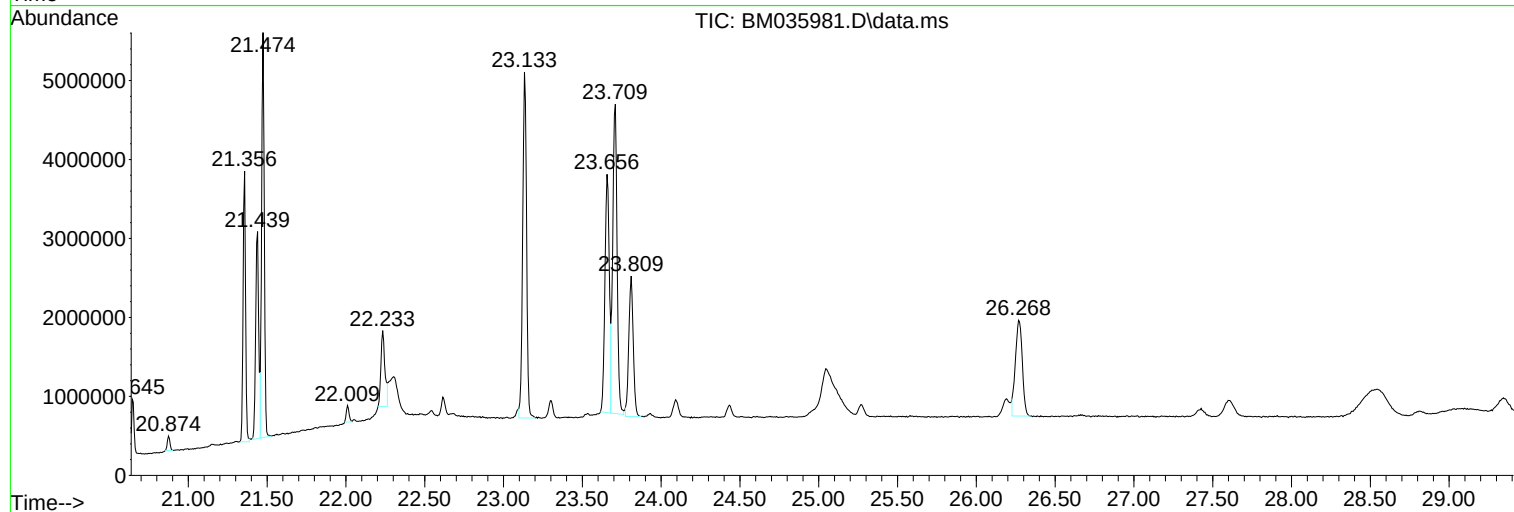
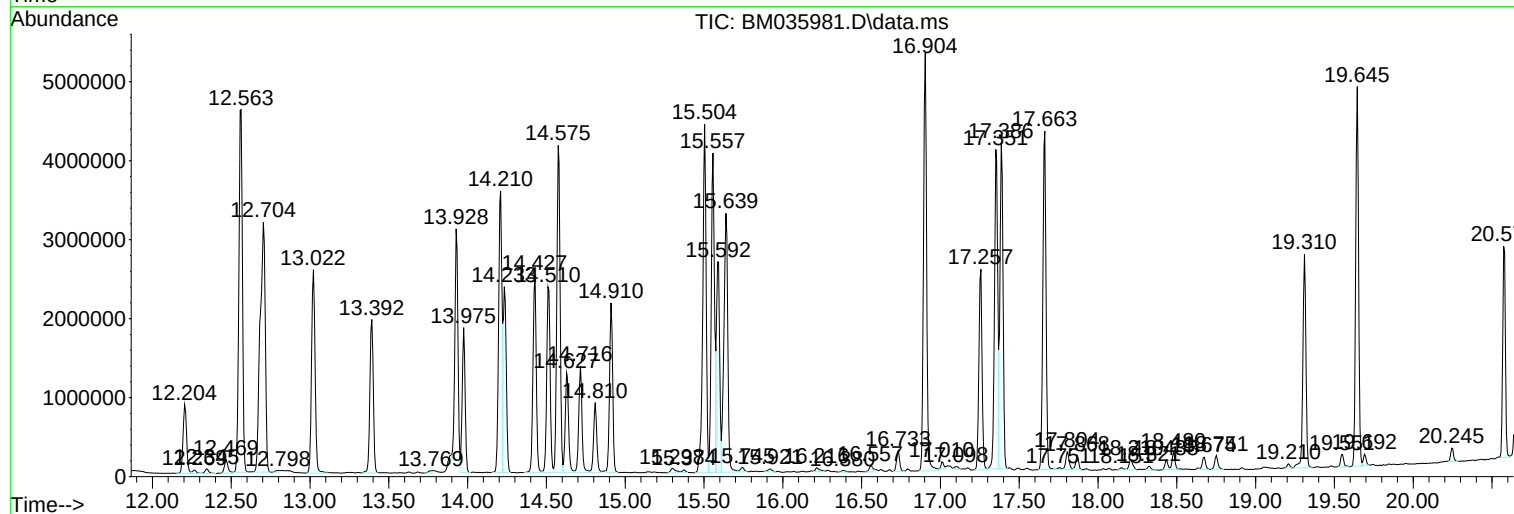
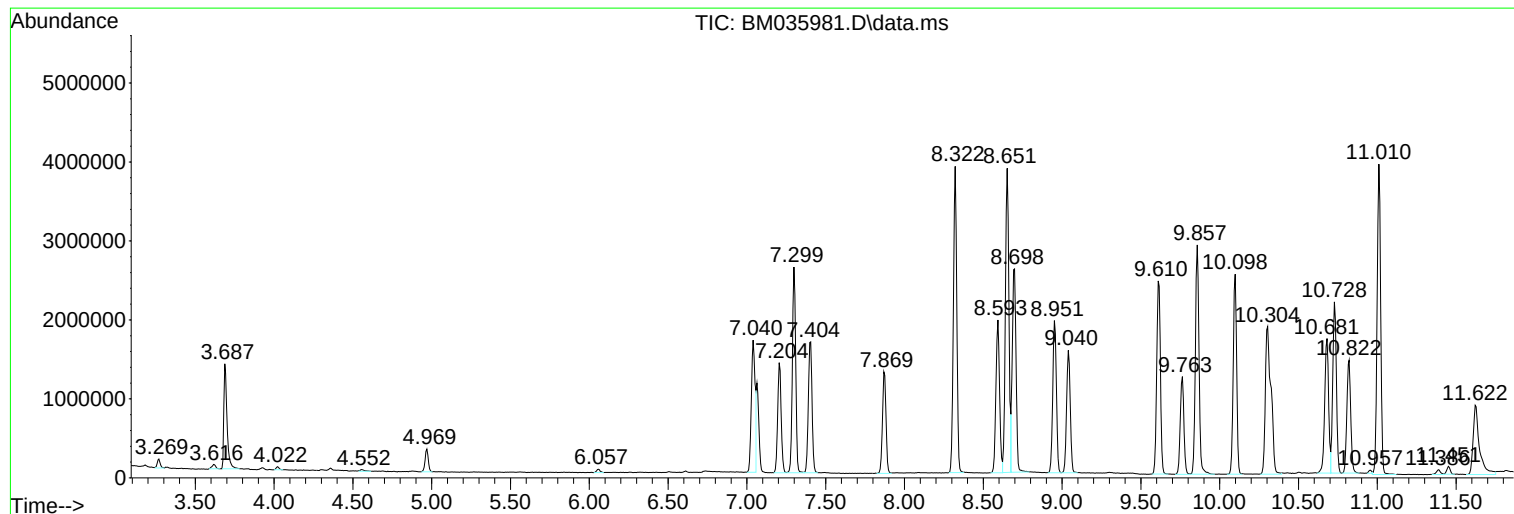
Sum of corrected areas: 247893390

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION

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



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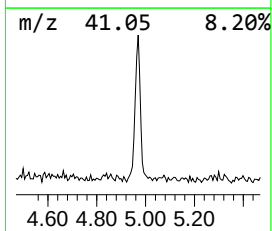
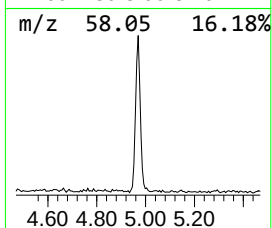
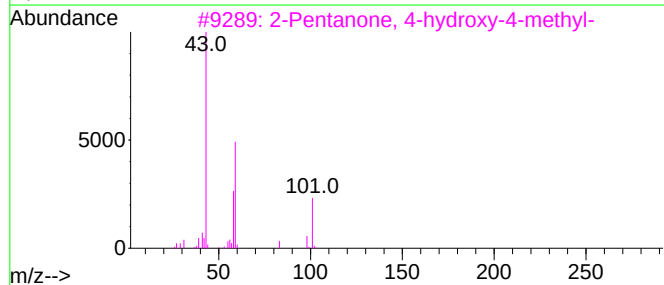
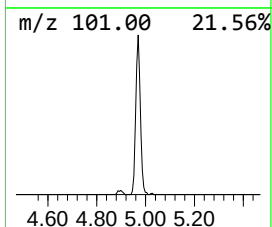
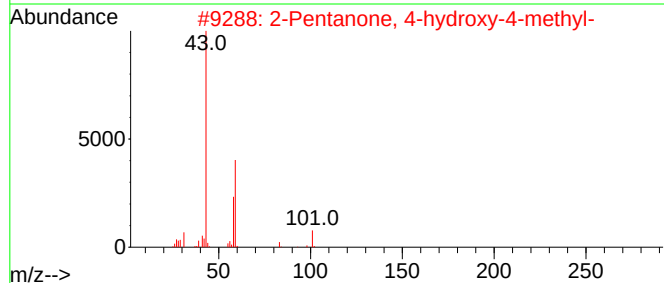
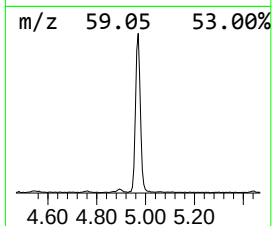
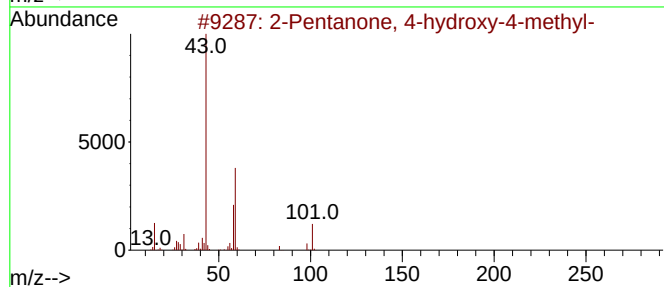
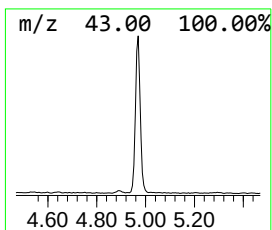
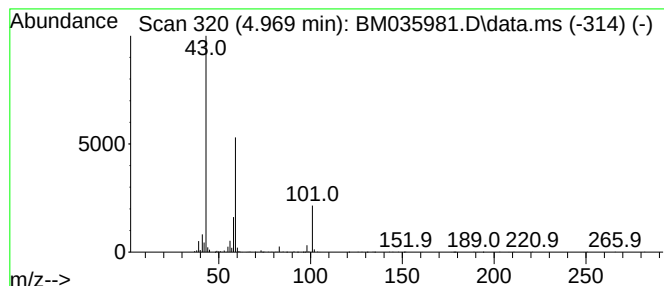
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.969	3.98 ng/ul	402353	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33		



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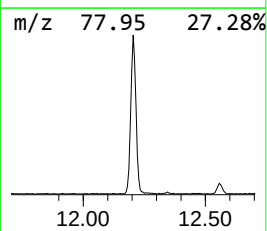
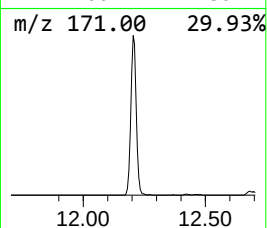
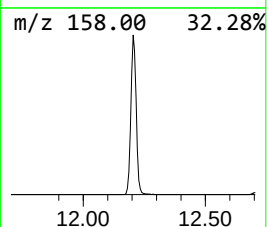
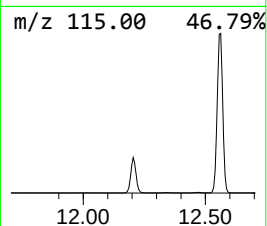
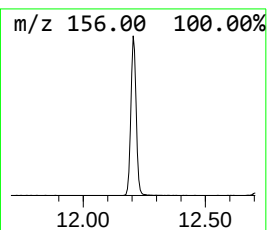
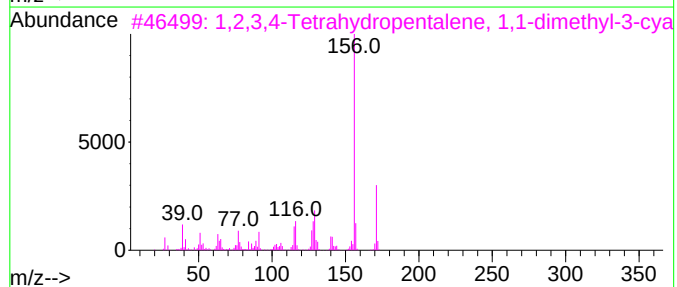
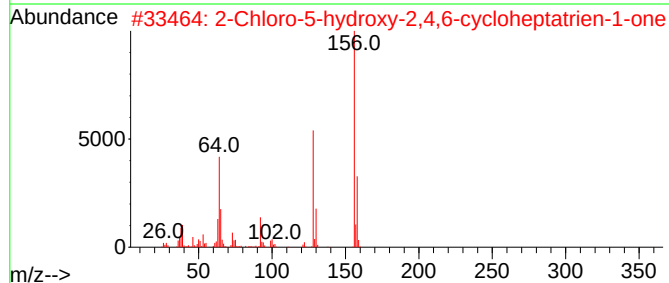
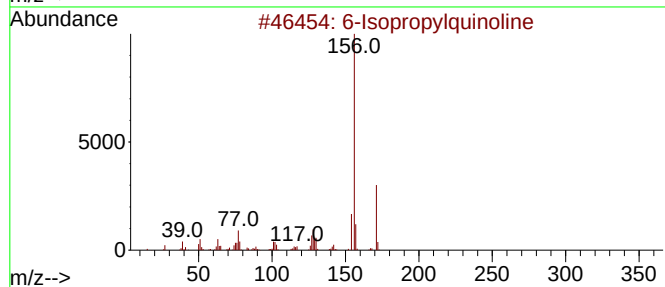
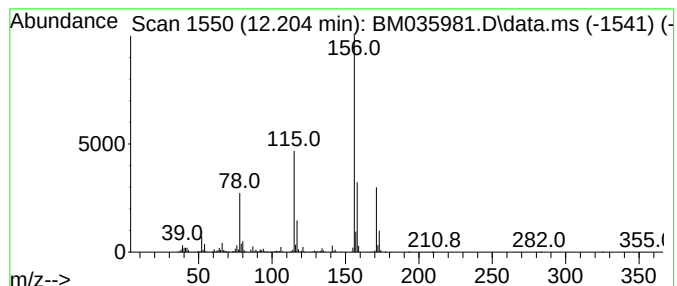
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.204	9.01 ng/ul	1335670	Naphthalene-d8	10.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Isopropylquinoline	171	C12H13N	000135-79-5	38		
2	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	38		
3	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37		
4	Chloroxylonol	156	C8H9ClO	000088-04-0	32		
5	Cyclopentanone, O-(trimethylsilyl...	171	C8H17NOSi	026208-87-7	32		



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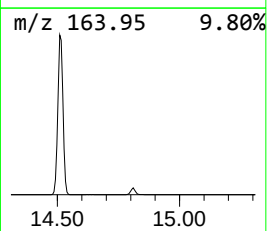
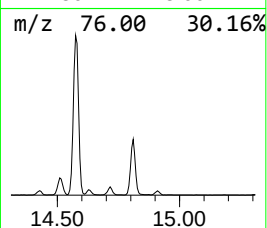
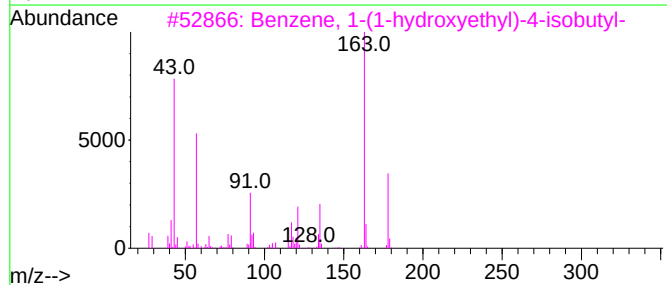
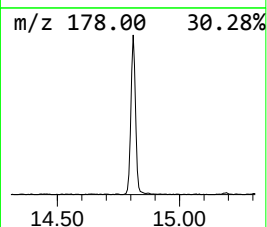
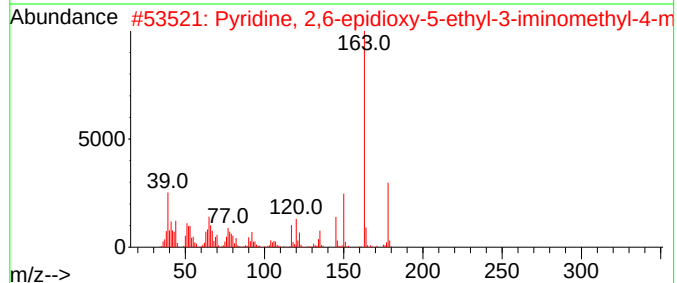
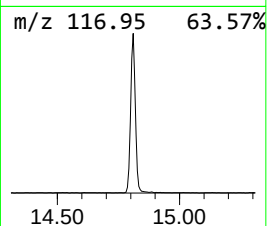
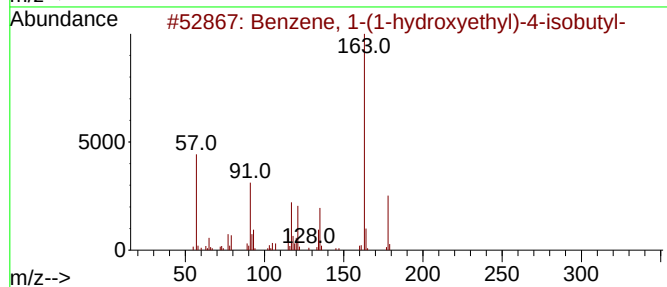
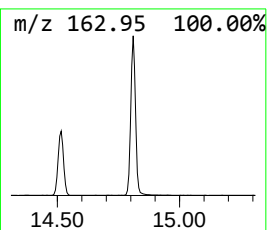
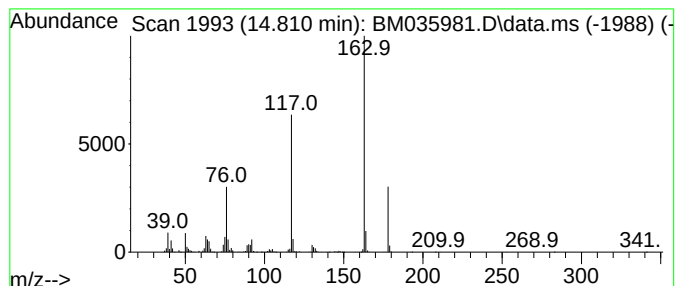
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1-(1-hydroxyethyl)... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.810	6.95 ng/ul	1235390	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	59
2			Pyridine, 2,6-epidioxo-5-ethyl-3...	178	C9H10N2O2	1000263-31-1	50
3			Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	50
4			2,3,4,5,6-Pentamethylphenol, met...	178	C12H18O	014804-37-6	47
5			Propofol	178	C12H18O	002078-54-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
Data File : BM035981.D
Acq On : 28 Jul 2022 12:17
Operator : CG/JU
Sample : N3874-08
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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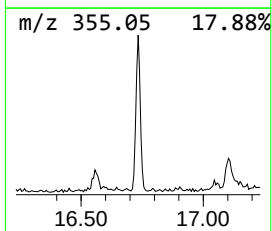
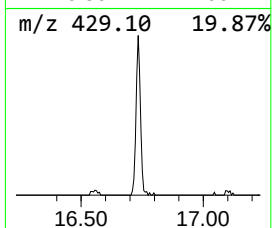
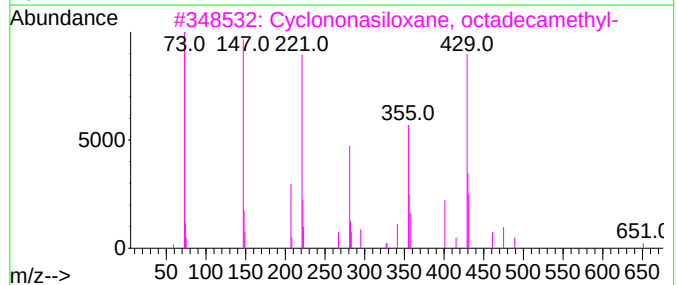
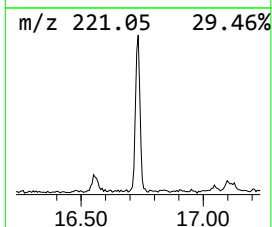
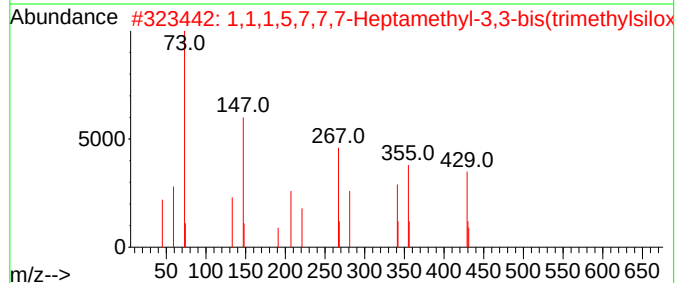
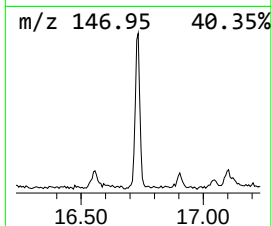
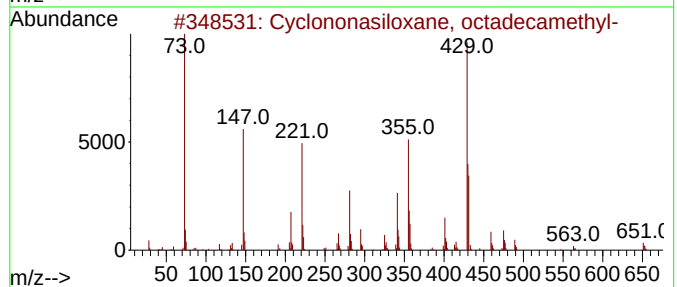
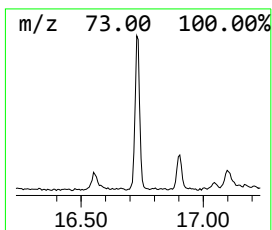
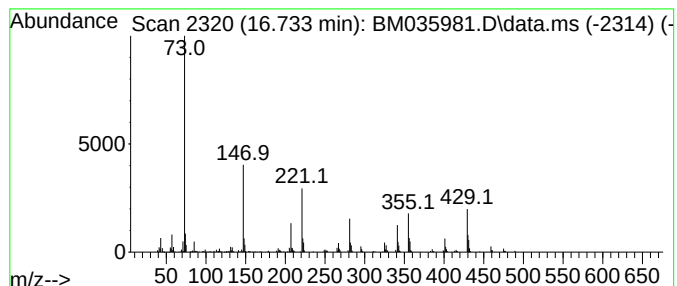
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclononasiloxane, octadeca... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.733	2.00 ng/ul	356915	Phenanthrene-d10	17.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	91
2			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	50
3			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	45
4			Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	38
5			Oxazepam, 2TMS derivative	430	C21H27ClN2O2Si2	055319-93-2	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
Data File : BM035981.D
Acq On : 28 Jul 2022 12:17
Operator : CG/JU
Sample : N3874-08
Misc :
ALS Vial : 38 Sample Multiplier: 1

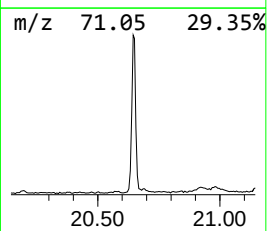
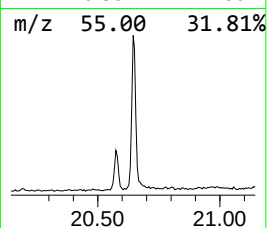
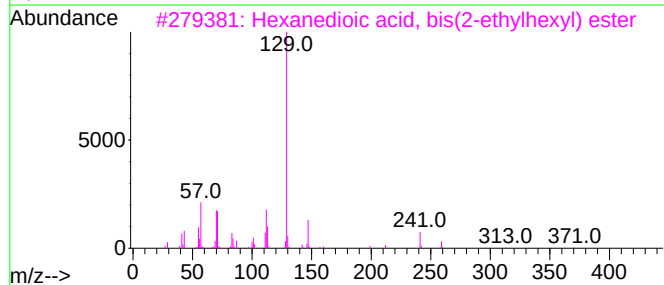
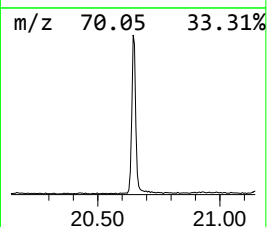
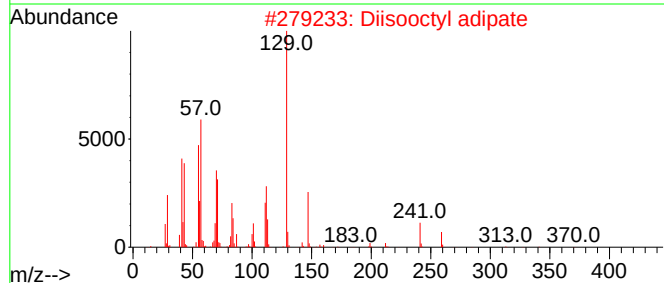
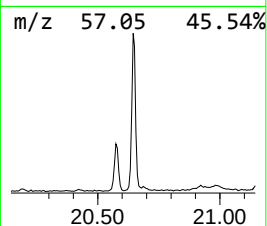
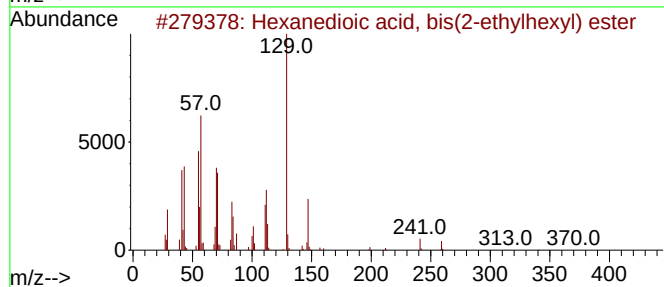
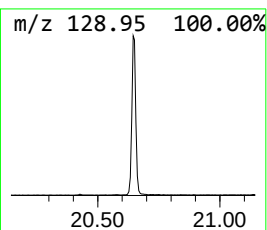
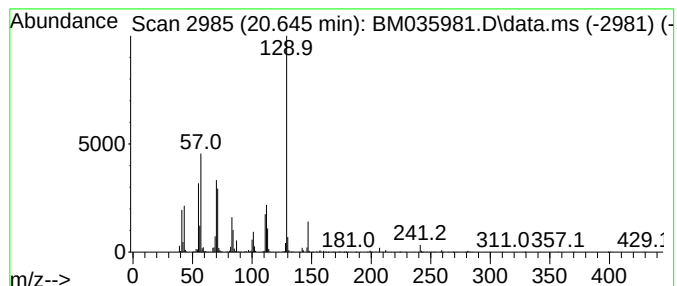
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexanedioic acid, bis(2-eth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.645	4.23 ng/ul	787530	Chrysene-d12	21.439	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	95
2	Diisooctyl adipate	370	C22H42O4	001330-86-5	94
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	74
5	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
Data File : BM035981.D
Acq On : 28 Jul 2022 12:17
Operator : CG/JU
Sample : N3874-08
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBJ34

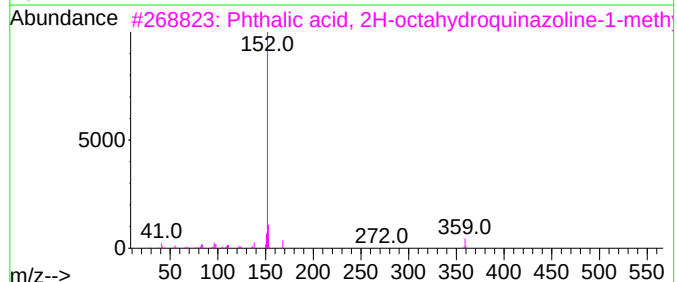
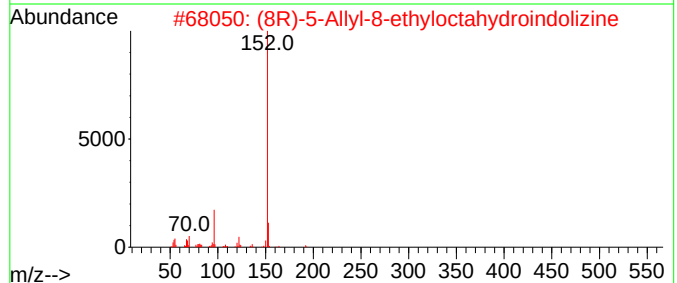
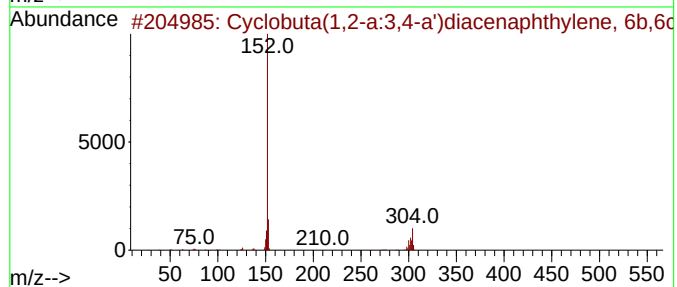
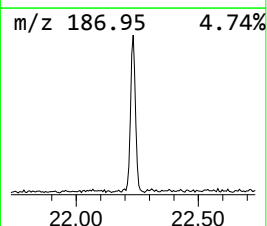
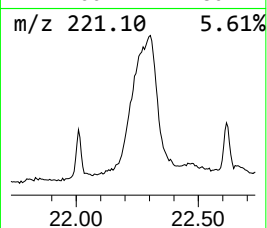
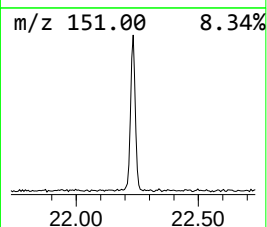
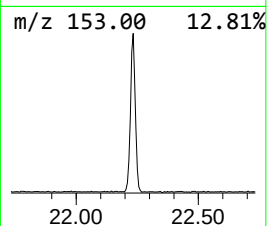
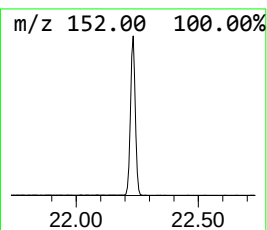
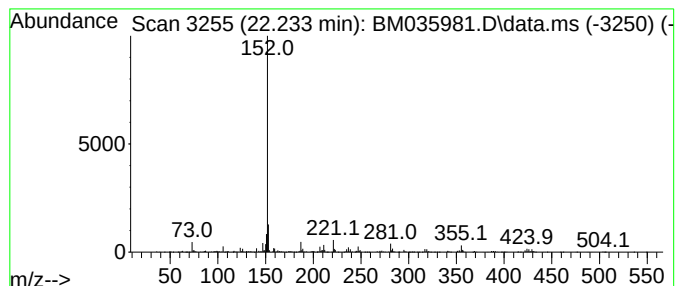
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cyclobuta(1,2-a:3,4-a')diac... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.233	9.28 ng/ul	1727650	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	64
2			(8R)-5-Allyl-8-ethyloctahydroind...	193	C13H23N	875587-44-3	64
3			Phthalic acid, 2H-octahydroquina...	359	C21H29NO4	1000322-80-2	64
4			2-Hydroxy-5-methoxybenzaldehyde,...	222	C12H14O4	1000447-33-2	50
5			Phthalic acid, heptyl 2H-octahyd...	415	C25H37NO4	1000322-80-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
Data File : BM035981.D
Acq On : 28 Jul 2022 12:17
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Sample : N3874-08
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
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
2-Pentanone, 4-...	4.969	4.0	ng/ul	402353	1	7.869	2021720	20.0
unknown-01	12.204	9.0	ng/ul	1335670	2	10.681	2963420	20.0
Benzene, 1-(1-h...	14.810	7.0	ng/ul	1235390	3	14.510	3556470	20.0
Cyclononasiloxa...	16.733	2.0	ng/ul	356915	4	17.257	3564300	20.0
Hexanedioic aci...	20.645	4.2	ng/ul	787530	5	21.439	3723380	20.0
Cyclobuta(1,2-a...	22.233	9.3	ng/ul	1727650	5	21.439	3723380	20.0