

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081524\
 Data File : BF139034.D
 Acq On : 15 Aug 2024 21:14
 Operator : RC/JU
 Sample : P3475-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-866-N-S0-0-0.5-080524

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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_F\Methods\8270-BF073024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139034.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	215	220	250	rBV	6093	31179	6.26%	0.743%
2	3.893	300	306	315	rVB	6642	12065	2.42%	0.288%
3	5.051	498	503	518	rBV	30544	47757	9.59%	1.138%
4	5.469	569	574	596	rBV	292196	442524	88.86%	10.549%
5	6.481	741	746	769	rBV	339451	498026	100.00%	11.872%
6	6.687	774	781	790	rBV2	5491	11306	2.27%	0.270%
7	6.834	801	806	815	rVB	205200	252002	50.60%	6.007%
8	7.398	897	902	922	rBV	218844	296369	59.51%	7.065%
9	8.116	1019	1024	1037	rBV	259607	326564	65.57%	7.785%
10	8.433	1074	1078	1097	rBV	24475	43082	8.65%	1.027%
11	9.186	1201	1206	1221	rBB	343929	427143	85.77%	10.182%
12	9.863	1316	1321	1332	rVB	280681	357370	71.76%	8.519%
13	9.957	1332	1337	1342	rBV	18869	27307	5.48%	0.651%
14	10.657	1451	1456	1468	rBV	210982	278390	55.90%	6.636%
15	11.351	1569	1574	1586	rBV	250788	332771	66.82%	7.933%
16	11.874	1659	1663	1678	rVB3	2580	6799	1.37%	0.162%
17	12.574	1777	1782	1786	rBV	12012	14830	2.98%	0.354%
18	12.804	1815	1821	1830	rBV	8820	14864	2.98%	0.354%
19	12.933	1838	1843	1848	rBV	245822	298180	59.87%	7.108%
20	13.833	1988	1996	2007	rBV7	5874	18987	3.81%	0.453%
21	13.998	2019	2024	2034	rVB	144675	198095	39.78%	4.722%
22	14.739	2145	2150	2156	rBV5	5382	8802	1.77%	0.210%
23	15.045	2197	2202	2204	rBV	6637	9951	2.00%	0.237%
24	15.074	2204	2207	2211	rVB3	5969	8090	1.62%	0.193%
25	15.409	2260	2264	2267	rVV	4086	5352	1.07%	0.128%
26	15.462	2267	2273	2287	rVB	126392	200571	40.27%	4.781%
27	15.868	2337	2342	2348	rBV4	6292	11306	2.27%	0.270%
28	16.368	2422	2427	2431	rBV4	2828	5165	1.04%	0.123%
29	16.945	2520	2525	2535	rBV7	3509	10155	2.04%	0.242%

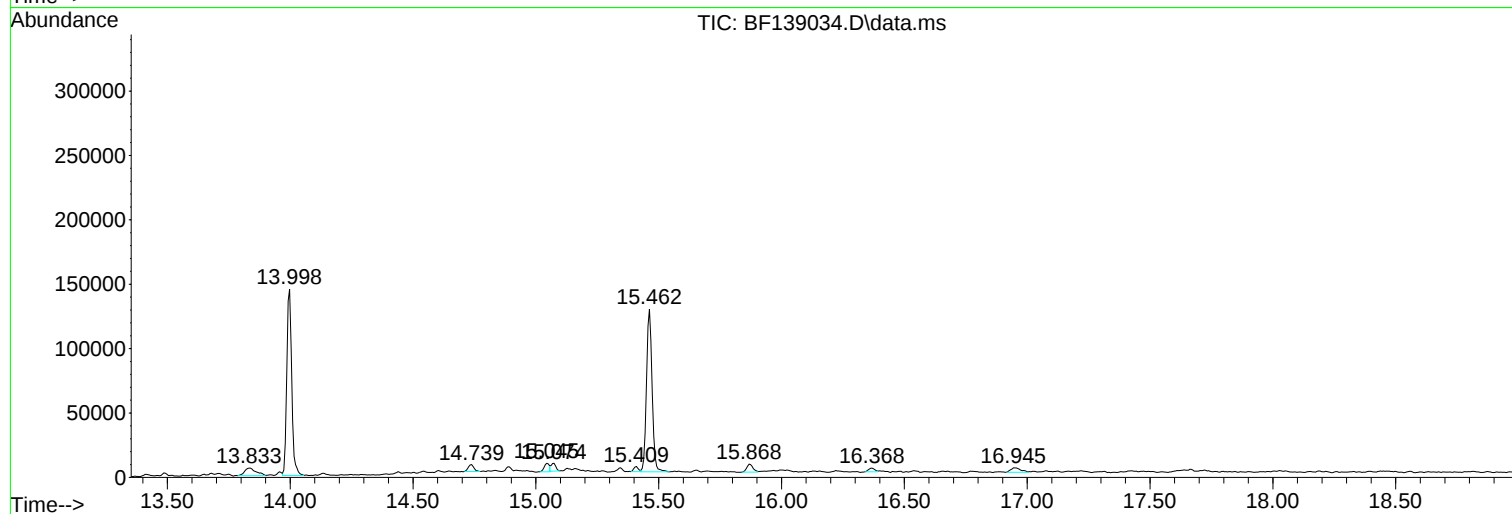
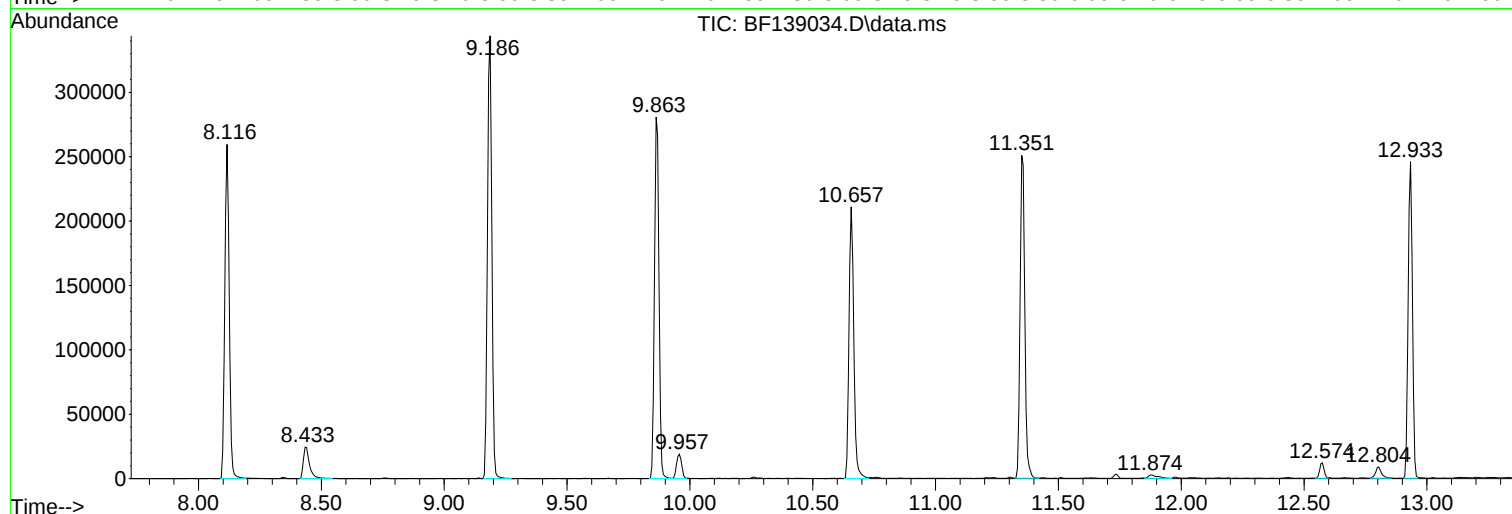
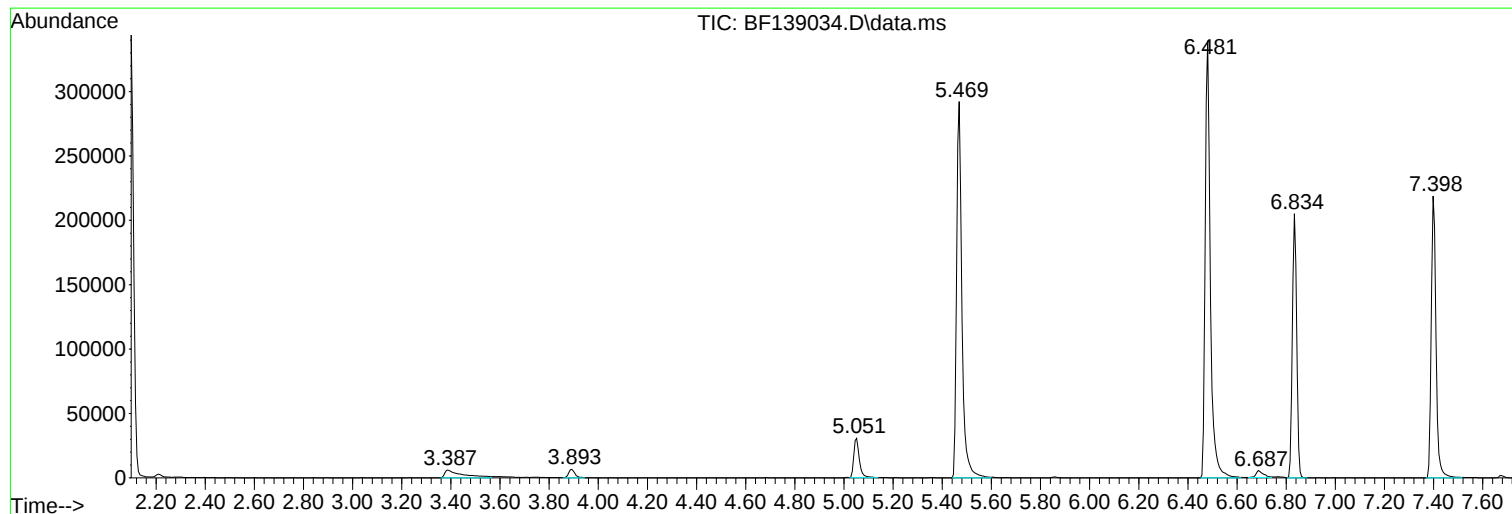
Sum of corrected areas: 4195002

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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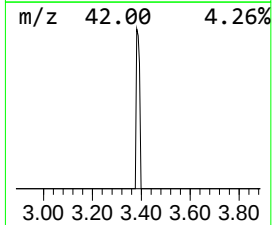
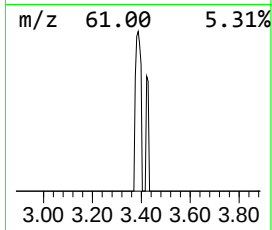
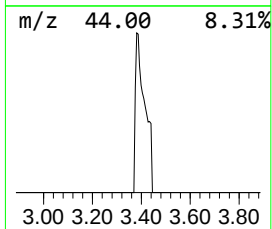
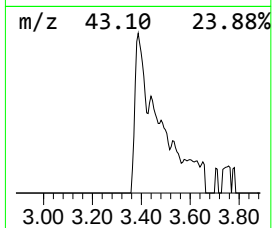
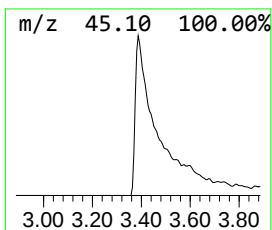
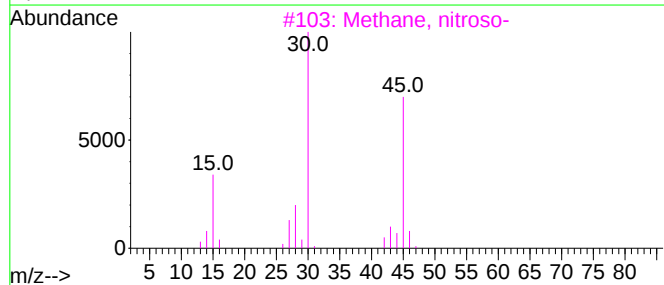
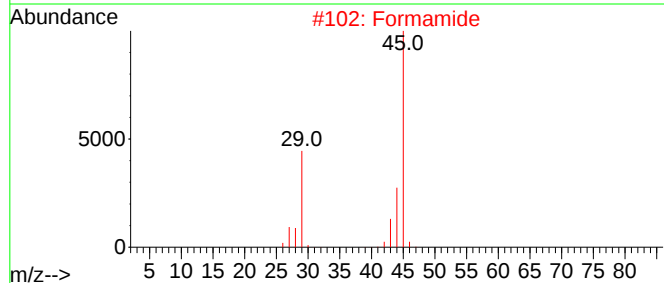
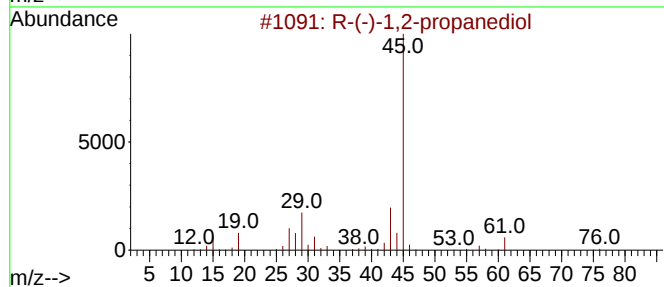
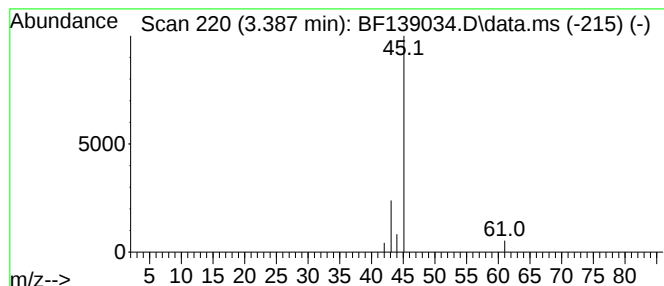
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.387 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.387	2.47 ng	31179	1,4-Dichlorobenzene-d4	6.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	4
2	Formamide			45	CH3NO	000075-12-7	3
3	Methane, nitroso-			45	CH3NO	000865-40-7	3
4	Propylene Glycol			76	C3H8O2	000057-55-6	2
5	Isopropyl Alcohol			60	C3H8O	000067-63-0	2



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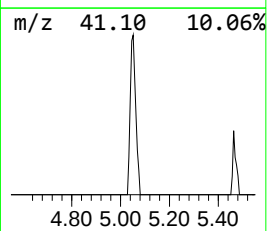
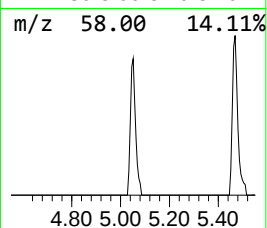
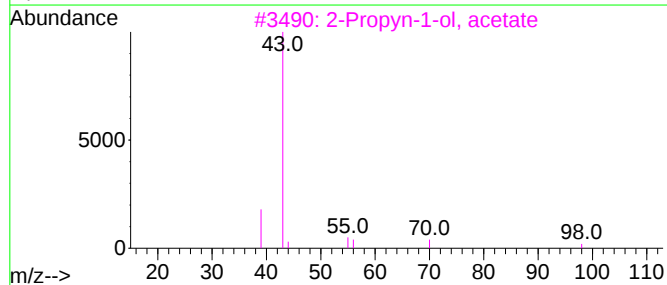
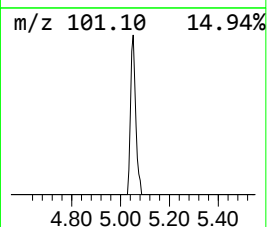
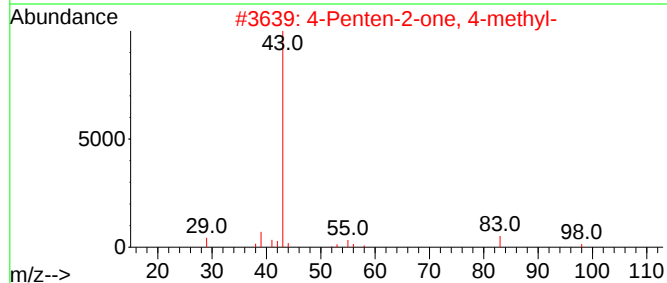
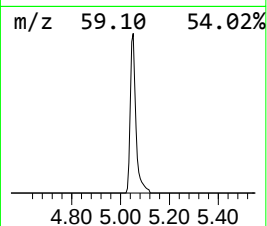
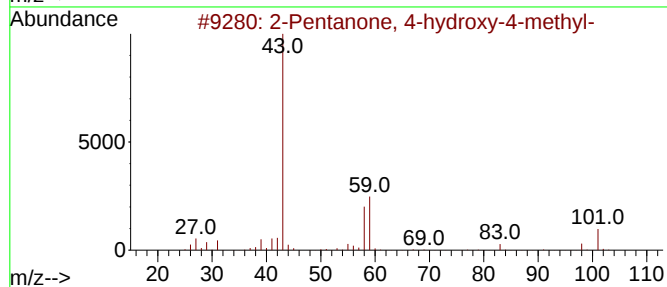
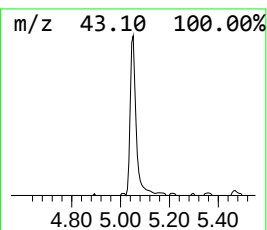
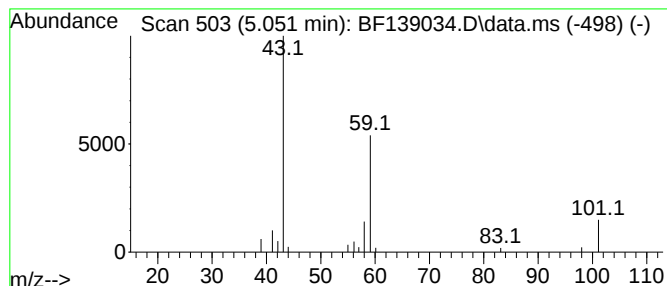
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.051	3.79 ng	47757	1,4-Dichlorobenzene-d4	6.834

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
3		2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
4		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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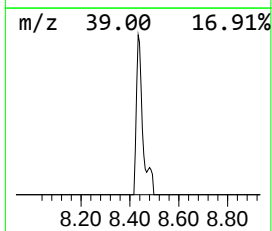
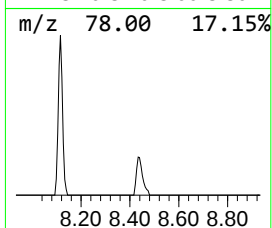
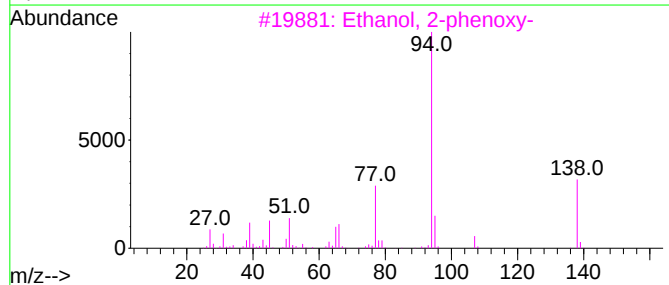
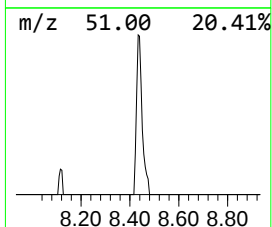
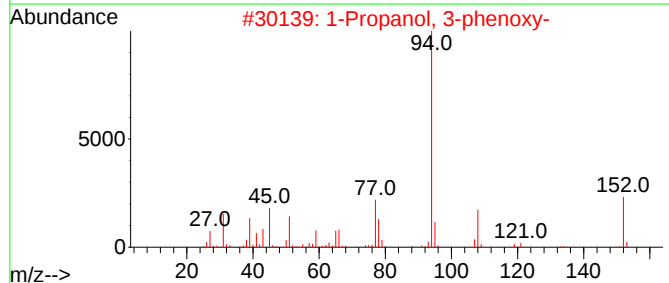
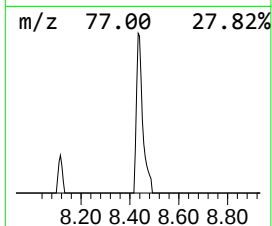
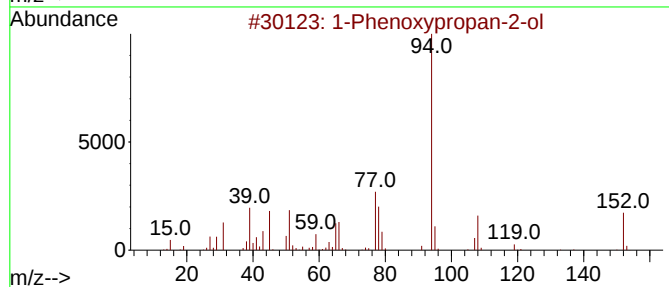
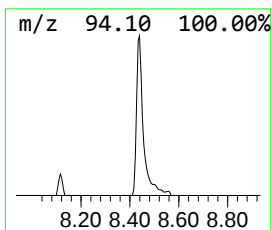
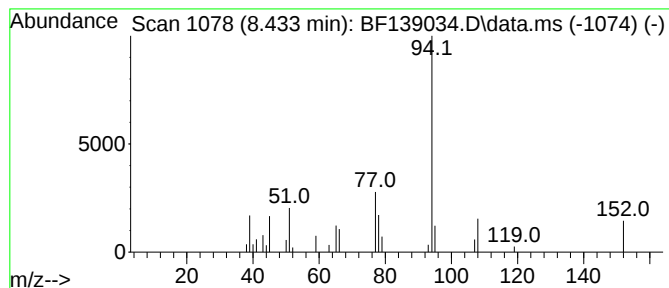
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.433	2.64 ng	43082	Naphthalene-d8	8.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
3			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	59
5			2-Hexen-4-yne, 2-methyl-	94	C7H10	058275-93-7	45



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.387	3.387	2.5	ng	31179	1	6.834	252002	20.0
2-Pentanone, 4-...	5.051	3.8	ng	47757	1	6.834	252002	20.0
1-Phenoxypropan...	8.433	2.6	ng	43082	2	8.116	326564	20.0